Access DB# 73851

SEARCH REQUEST FORM

Scientific and Technical Information Center

Requester's Full Name: Hong Liu Examiner #: _____ Date: 8/20/02

Art Unit: 1624 Phone Number 30 6-5814 Serial Number: 09/864,905

Mail Box and Bldg/Room Location: 4E01 Results Format Preferred (circle): PAPER DISK E-MAIL

If more than one search is submitted, please prioritize searches in order of need.

****

Please provide a detailed statement of the search topic, and describe as specifically as possible the subject matter to be searched. Include the elected species or structures, keywords, synonyms, acronyms, and registry numbers, and combine with the concept or utility of the invention. Define any terms that may have a special meaning. Give examples or relevant citations, authors, etc, if known. Please attach a copy of the cover sheet, pertinent claims, and abstract.

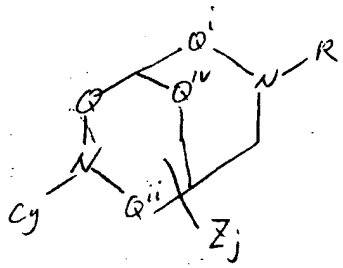
Title of Invention: Pharm. compositions & methods for use

Inventors (please provide full names): Miller, C Dull, G Miao, L Lynn, D
Schmitt, J Clark, T

Earliest Priority Filing Date: _____

***For Sequence Searches Only* Please include all pertinent information (parent, child, divisional, or issued patent numbers) along with the appropriate serial number.**

Barb please!



Point of Contact:
Barb O'Bryen
Technical Information Specialist
STIC CM1 6A05 308-4291

Applicants elected the species of 3-(5-phenoxy-3-pyridyl)-3,7-diazabicyclo[3,3,1]nonane. Please based the structure search on this core.

****

STAFF USE ONLY	Type of Search	Vendors and cost where applicable
Searcher: BOB	NA Sequence (#) _____	STN 242
Searcher Phone #: _____	AA Sequence (#) _____	Dialog _____
Searcher Location: _____	Structure (#) 3	Questel/Orbit _____
Date Searcher Picked Up: _____	Bibliographic _____	Dr.Link _____
Date Completed: 8-30-02	Litigation _____	Lexis/Nexis _____
Searcher Prep & Review Time: 40	Fulltext _____	Sequence Systems _____
Clerical Prep Time: _____	Patent Family _____	WWW/Internet _____
Online Time: 35	Other _____	Other (specify) Chem Draw

PTO-1590 (8-01)

species for wo 97/40049

Welcome to STN International! Enter x:x

LOGINID:ssspta1611hxl

PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * * * * * * Welcome to STN International * * * * * * * * * *

NEWS 1		Web Page URLs for STN Seminar Schedule - N. America
NEWS 2	Apr 08	"Ask CAS" for self-help around the clock
NEWS 3	Apr 09	BEILSTEIN: Reload and Implementation of a New Subject Area
NEWS 4	Apr 09	ZDB will be removed from STN
NEWS 5	Apr 19	US Patent Applications available in IFICDB, IFIPAT, and IFIUDB
NEWS 6	Apr 22	Records from IP.com available in CAPLUS, HCAPLUS, and ZCAPLUS
NEWS 7	Apr 22	BIOSIS Gene Names now available in TOXCENTER
NEWS 8	Apr 22	Federal Research in Progress (FEDRIP) now available
NEWS 9	Jun 03	New e-mail delivery for search results now available
NEWS 10	Jun 10	MEDLINE Reload
NEWS 11	Jun 10	PCTFULL has been reloaded
NEWS 12	Jul 02	FOREGE no longer contains STANDARDS file segment
NEWS 13	Jul 22	USAN to be reloaded July 28, 2002; saved answer sets no longer valid
NEWS 14	Jul 29	Enhanced polymer searching in REGISTRY
NEWS 15	Jul 30	NETFIRST to be removed from STN
NEWS 16	Aug 08	CANCERLIT reload
NEWS 17	Aug 08	PHARMAMarketLetter(PHARMAML) - new on STN
NEWS 18	Aug 08	NTIS has been reloaded and enhanced
NEWS 19	Aug 19	Aquatic Toxicity Information Retrieval (AQUIRE) now available on STN
NEWS 20	Aug 19	IFIPAT, IFICDB, and IFIUDB have been reloaded
NEWS 21	Aug 19	The MEDLINE file segment of TOXCENTER has been reloaded
NEWS 22	Aug 26	Sequence searching in REGISTRY enhanced
NEWS 23	Sep 03	JAPIO has been reloaded and enhanced
NEWS 24	Sep 16	Experimental properties added to the REGISTRY file
NEWS 25	Sep 16	Indexing added to some pre-1967 records in CA/CAPLUS
NEWS 26	Sep 16	CA Section Thesaurus available in CAPLUS and CA
NEWS 27	Oct 01	CASREACT Enriched with Reactions from 1907 to 1985

NEWS EXPRESS	February 1 CURRENT WINDOWS VERSION IS V6.0d, CURRENT MACINTOSH VERSION IS V6.0a(ENG) AND V6.0Ja(JP), AND CURRENT DISCOVER FILE IS DATED 05 FEBRUARY 2002
NEWS HOURS	STN Operating Hours Plus Help Desk Availability
NEWS INTER	General Internet Information
NEWS LOGIN	Welcome Banner and News Items
NEWS PHONE	Direct Dial and Telecommunication Network Access to STN
NEWS WWW	CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that

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* * * * * * * * * * * * * * STN Columbus * * * * * * * * * * * * * * *

FILE 'HOME' ENTERED AT 17:57:47 ON 04 OCT 2002

=> fil reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 17:57:55 ON 04 OCT 2002
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 3 OCT 2002 HIGHEST RN 459123-02-5
DICTIONARY FILE UPDATES: 3 OCT 2002 HIGHEST RN 459123-02-5

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:
http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

=> list saved
LIST IS NOT A RECOGNIZED COMMAND
The previous command name entered was not recognized by the system.
For a list of commands available to you in the current file, enter
"HELP COMMANDS" at an arrow prompt (=>).

=> saved
SAVED IS NOT A RECOGNIZED COMMAND
The previous command name entered was not recognized by the system.
For a list of commands available to you in the current file, enter
"HELP COMMANDS" at an arrow prompt (=>).

```
=> help commands
Enter one of these commands at the arrow prompt (=>).

ACTIVATE ----- Assign L#s to saved query or answer set.
ANALYZE ------ Build expand terms from answer fields.
ARCHIVE ------ Purchase rights for archiving.
BATCH -------- Request a batch search.
DELETE ------- Delete saved or current session items.
DISPLAY ------ Display saved or current session items.
DOWNLOAD ----- Transfer data using Kermit protocol.
EDIT --------- Modify the text of an E-number entry.
EXPAND ------- Look at the index around a term.
FILE --------- Specify the search and display file.
HELP --------- For help on how to use the system.
INDEX -------- Specify the Index environment.
LOGOFF ------- End the online session.
NEWS --------- Display current news about the system.
ORDER -------- Order an original document or copy.
POLYLINK ----- Search Using Polymer Linking
PRINT -------- Print answers offline.
QUERY -------- Define a search question (query).
REDISTRIBUTE - Purchase rights for redistribution.
SAVE --------- Save an L-numbered query or answer set.
SCREEN ------- Define a set of screens for searching.
SDI ---------- Request searches be run on file updates.
SEARCH ------- Perform a search.
SELECT ------- Build expand terms from answer fields.
SEQLINK ------ Search Using Sequence Linking
SET ---------- Set terminal and interaction options.
SORT --------- Sort Answers of an L-number list
STRUCTURE ---- Create a structure for searching.
TABULATE ----- Analyze SmartSELECT L-number data.
TRANSFER ----- Extract terms from answers and search.
? ------------ The same as HELP.
Enter one of these commands at the arrow prompt (=>).


For more detailed information about a command, enter HELP followed by
the command name at an arrow prompt.

Example: => HELP SEARCH

For information on entering multiple commands at a single arrow
prompt, enter HELP COMMAND STACKING.

=> display saved
NAME             CREATED      NOTES/TITLE
--------------   -----------  -----------------------------------
F403066/A        07 JUL 2000  3466 ANSWERS IN FILE REGISTRY
F485188/A        22 AUG 2000  15 ANSWERS IN FILE REGISTRY
F582302/A        29 MAR 2001  49 ANSWERS IN FILE REGISTRY
F9367871/A       15 MAY 2002  1856 ANSWERS IN FILE REGISTRY
F9386258/A       19 SEP 2000  179 ANSWERS IN FILE REGISTRY
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F9396156/A       05 JUL 2000  28 ANSWERS IN FILE REGISTRY
F9445504/A       02 JUN 2000  63 ANSWERS IN FILE REGISTRY
F9446147/A       14 AUG 2000  301 ANSWERS IN FILE REGISTRY
F9446423/A       27 JUN 2000  107 ANSWERS IN FILE REGISTRY
F9462797/A       16 JUL 2000  13 ANSWERS IN FILE REGISTRY
F9462812/A       19 SEP 2000  0 ANSWERS IN FILE REGISTRY
F9482810/A       15 AUG 2000  226 ANSWERS IN FILE REGISTRY
F9486327/A       15 JUL 2000  51 ANSWERS IN FILE REGISTRY
F9486646/A       08 MAR 2001  12669 ANSWERS IN FILE REGISTRY
F9489271/A       02 JAN 2001  155 ANSWERS IN FILE REGISTRY
F9499699/A       01 JUN 2000  7 ANSWERS IN FILE REGISTRY
F9501758/A       11 OCT 2000  42 ANSWERS IN FILE REGISTRY
F9504183/A       26 JUN 2000  3 ANSWERS IN FILE REGISTRY
F9508240/A       16 DEC 2000  363 ANSWERS IN FILE REGISTRY
F9508796/A       12 JUL 2000  91 ANSWERS IN FILE REGISTRY
F9508989/A       27 FEB 2001  176 ANSWERS IN FILE REGISTRY
F9509831/A       07 SEP 2000  165 ANSWERS IN FILE REGISTRY
F9513147/A       15 JUN 2000  360 ANSWERS IN FILE REGISTRY
F9538800/A       29 DEC 2000  114 ANSWERS IN FILE REGISTRY
F954237/A        26 SEP 2000  7030 ANSWERS IN FILE REGISTRY
F9546555/A       21 JUL 2000  2536 ANSWERS IN FILE REGISTRY
F9552036/A       21 FEB 2001  116 ANSWERS IN FILE REGISTRY
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F9566002/A       25 JUL 2000  201 ANSWERS IN FILE REGISTRY
F9583924/A       23 DEC 2000  910 ANSWERS IN FILE REGISTRY
F9592336/A       23 AUG 2000  141 ANSWERS IN FILE REGISTRY
F9606921/A       25 OCT 2000  167 ANSWERS IN FILE REGISTRY
F9622086/A       10 JAN 2001  136 ANSWERS IN FILE REGISTRY
F9627766/A       28 NOV 2000  0 ANSWERS IN FILE REGISTRY
F9634801/A       14 OCT 2000  417 ANSWERS IN FILE REGISTRY
F9687241/A       08 JUN 2001  195 ANSWERS IN FILE REGISTRY
F9693896/A       31 JAN 2001  9183 ANSWERS IN FILE REGISTRY
F9696504/A       15 MAY 2001  4549 ANSWERS IN FILE REGISTRY
F9718223/A       25 MAY 2001  226 ANSWERS IN FILE REGISTRY
F9787672/L       14 JUN 2001  2 L-NUMBERS
F9792866/A       25 SEP 2001  1833 ANSWERS IN FILE REGISTRY
F9868992/A       25 JAN 2002  574 ANSWERS IN FILE REGISTRY
F9977944/A       17 JAN 2002  771 ANSWERS IN FILE REGISTRY
R9380742/A       23 JUN 2000  114 ANSWERS IN FILE REGISTRY
R9655115/A       01 MAR 2001  322 ANSWERS IN FILE REGISTRY


=> delete saved
DELETE ALL SAVED QUERIES, STRUCTURES, SCREEN SETS, ANSWER SETS AND L# LISTS?
(Y)/N:y
ALL SAVED SCREENS, STRUCTURES, QUERIES, ANSWER SETS, AND L# LISTS DELETED

=> fil caplus
COST IN U.S. DOLLARS                            SINCE FILE      TOTAL
                                                   ENTRY    SESSION
FULL ESTIMATED COST                                 1.14       1.35

FILE 'CAPLUS' ENTERED AT 17:59:49 ON 04 OCT 2002
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
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PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

FILE COVERS 1907 - 4 Oct 2002 VOL 137 ISS 15
FILE LAST UPDATED: 3 Oct 2002 (20021003/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter HELP ROLES at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

=> s wo9740049/pn
L1 1 WO9740049/PN

=> select l1 rn 1-
E1 THROUGH E197 ASSIGNED

=> fil reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	2.11	3.46

FILE 'REGISTRY' ENTERED AT 18:00:27 ON 04 OCT 2002

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 3 OCT 2002 HIGHEST RN 459123-02-5
DICTIONARY FILE UPDATES: 3 OCT 2002 HIGHEST RN 459123-02-5

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:
http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

=> s e1-e197
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1 4584-46-7/BI
(4584-46-7/RN)
1 4755-77-5/BI
(4755-77-5/RN)
1 510-77-0/BI
(510-77-0/RN)
1 51186-58-4/BI
(51186-58-4/RN)
1 5241-66-7/BI
(5241-66-7/RN)
1 5472-49-1/BI
(5472-49-1/RN)
1 563-41-7/BI
(563-41-7/RN)
1 57-14-7/BI
(57-14-7/RN)
1 574-98-1/BI
(574-98-1/RN)
1 592-82-5/BI
(592-82-5/RN)
1 60-34-4/BI
(60-34-4/RN)
1 60384-53-4/BI
(60384-53-4/RN)

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              1 624-84-0/BI
                (624-84-0/RN)
              1 65232-62-4/BI
                (65232-62-4/RN)
              1 7250-67-1/BI
                (7250-67-1/RN)
              1 75-36-5/BI
                (75-36-5/RN)
              1 7536-58-5/BI
                (7536-58-5/RN)
              1 7803-57-8/BI
                (7803-57-8/RN)
              1 79-07-2/BI
                (79-07-2/RN)
              1 79-22-1/BI
                (79-22-1/RN)
              1 86-84-0/BI
                (86-84-0/RN)
              1 870-46-2/BI
                (870-46-2/RN)
              1 96763-06-3/BI
                (96763-06-3/RN)
L2          197 (100-39-0/BI OR 100944-15-8/BI OR 103-71-9/BI OR 103-72-0/BI
OR
                105-36-2/BI OR 105-39-5/BI OR 106-95-6/BI OR 106-96-7/BI OR
                107-13-1/BI OR 107-14-2/BI OR 109-70-6/BI OR 109-84-2/BI OR
                110-53-2/BI OR 111-25-1/BI OR 111-36-4/BI OR 112-67-4/BI OR
                122584-14-9/BI OR 125145-51-9/BI OR 13734-34-4/BI OR
14649-03-7/
                BI OR 1576-35-8/BI OR 179107-98-3/BI OR 179107-99-4/BI OR
179108
                -00-0/BI OR 179108-02-2/BI OR 179108-04-4/BI OR 179108-05-5/BI
                OR 179108-06-6/BI OR 179239-41-9/BI OR 180854-29-9/BI OR
183487-
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198987-82-
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                198987-86-9/BI OR 198987-87-0/BI OR 198987-88-1/BI OR
198987-89-
                2/BI OR 198987-90-5/BI OR 198987-91-6/BI OR 198987-92-7/BI OR
                198987-93-8/BI OR 198987-94-9/BI OR 198987-95-0/B

=> d scan

L2   197 ANSWERS   REGISTRY  COPYRIGHT 2002 ACS
IN   6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 11-ethyl-4a,5,9,10,11,12-
     hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF   C18 H23 N O3

Relative stereochemistry.
```
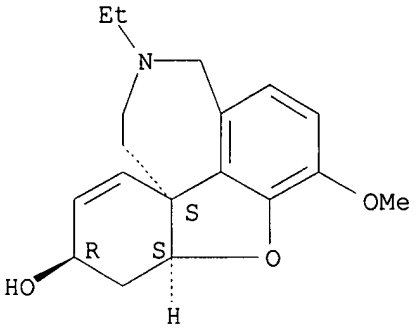
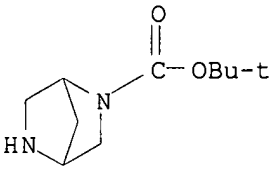



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

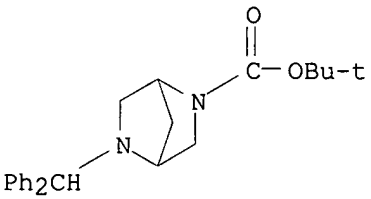
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):196

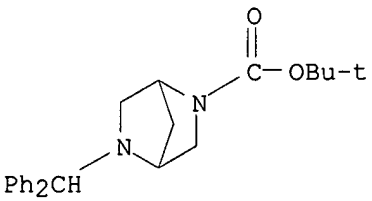
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane-2-carboxylic acid, 1,1-dimethylethyl ester (9CI)
MF C10 H18 N2 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

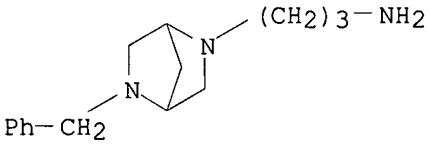
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane-2-carboxylic acid, 5-(diphenylmethyl)-, 1,1-dimethylethyl ester (9CI)
MF C23 H28 N2 O2





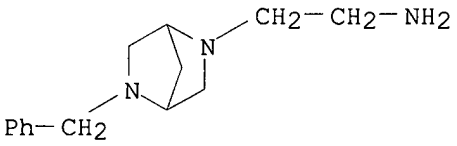
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane-2-propanamine, 5-(phenylmethyl)- (9CI)
MF C15 H23 N3



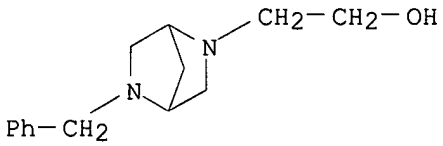
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane-2-ethanamine, 5-(phenylmethyl)- (9CI)
MF C14 H21 N3



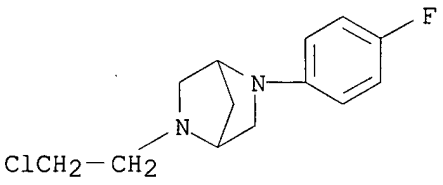
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane-2-ethanol, 5-(phenylmethyl)- (9CI)
MF C14 H20 N2 O



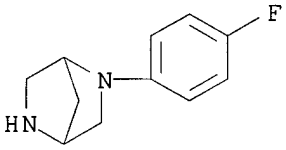
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(2-chloroethyl)-5-(4-fluorophenyl)- (9CI)
MF C13 H16 Cl F N2



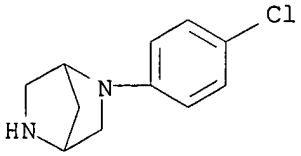
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-fluorophenyl)- (9CI)
MF C11 H13 F N2



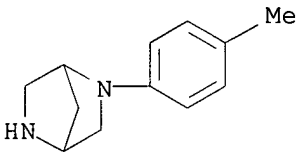
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-chlorophenyl)- (9CI)
MF C11 H13 Cl N2



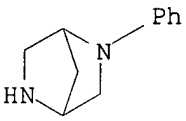
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-methylphenyl)- (9CI)
MF C12 H16 N2



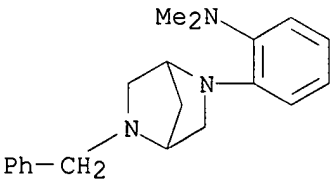
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-phenyl- (9CI)
MF C11 H14 N2
CI COM



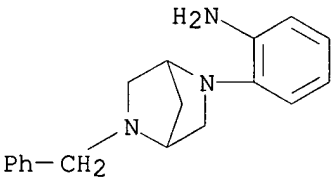
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenamine,
N,N-dimethyl-2-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-
2-yl]- (9CI)
MF C20 H25 N3



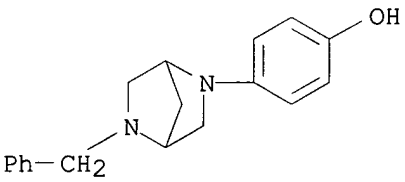
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenamine, 2-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]- (9CI)
MF C18 H21 N3



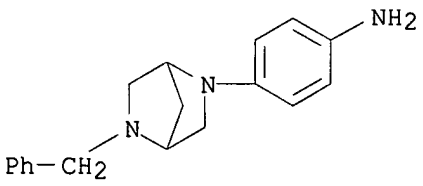
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Phenol, 4-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]- (9CI)
MF C18 H20 N2 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

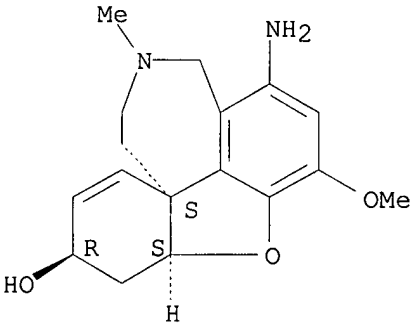
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenamine, 4-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]- (9CI)
MF C18 H21 N3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-amino-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aS,6R,8aS)- (9CI)
MF C17 H22 N2 O3

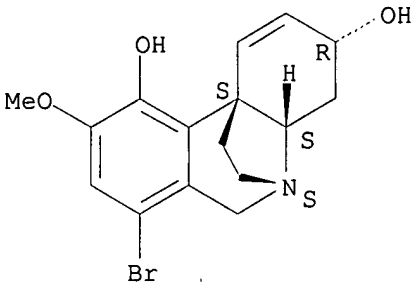
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 3H,6H-5,10b-Ethanophenanthridine-3,10-diol, 7-bromo-4,4a-dihydro-9-methoxy-, (3.alpha.,4a.beta.,5.beta.,10b.beta.)- (9CI)
MF C16 H18 Br N O3

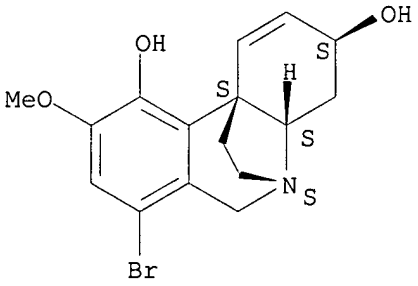
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 3H,6H-5,10b-Ethanophenanthridine-3,10-diol, 7-bromo-4,4a-dihydro-9-methoxy-, (3.alpha.,4a.alpha.,5.alpha.,10b.alpha.)- (9CI)
MF C16 H18 Br N O3

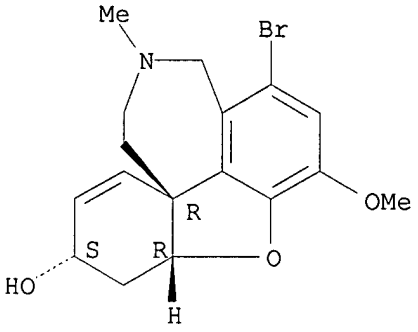
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, [4aR-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C17 H20 Br N O3

Absolute stereochemistry.

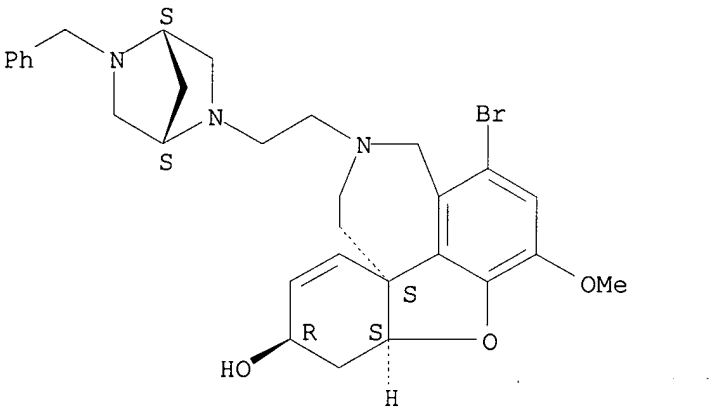


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-

hexahydro-3-methoxy-11-[2-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]ethyl]-, [4a.alpha.,6.beta.,8aR*,11(1R*,4R*)]- (9CI)
MF C30 H36 Br N3 O3

Relative stereochemistry.

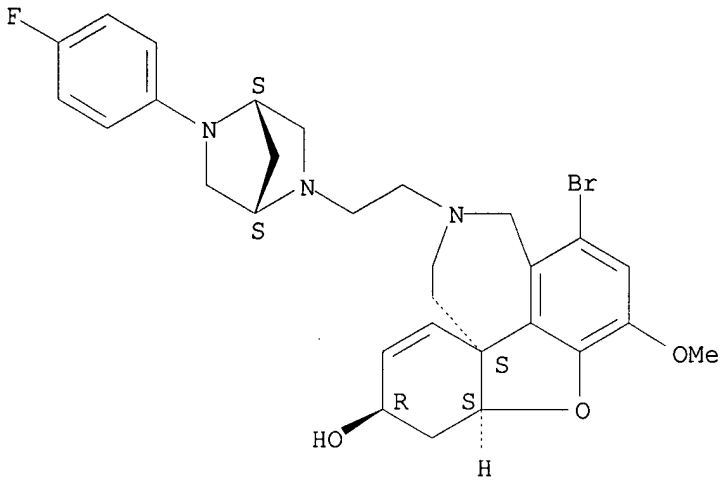


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-11-[2-[5-(4-fluorophenyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]ethyl]-4a,5,9,10,11,12-hexahydro-3-methoxy-, [4a.alpha.,6.beta.,8aR*,11(1R*,4R*)]- (9CI)

MF C29 H33 Br F N3 O3

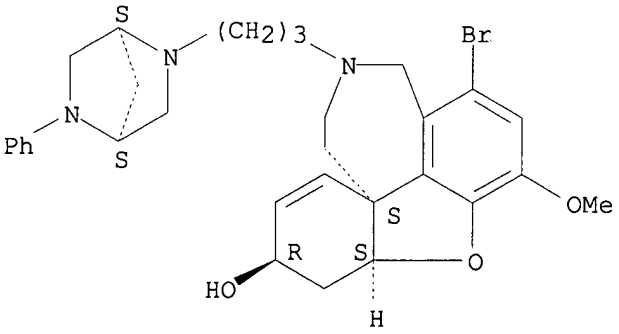
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-[3-(5-phenyl-2,5-diazabicyclo[2.2.1]hept-2-yl)propyl]-, [4a.alpha.,6.beta.,8aR*,11(1R*,4R*)]- (9CI)
MF C30 H36 Br N3 O3

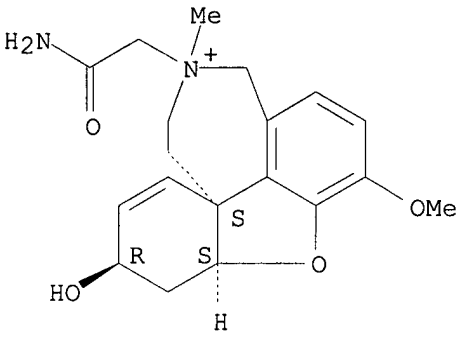
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 11-(2-amino-2-oxoethyl)-4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-, bromide, (4aS,6R,8aS)-[partial]- (9CI)
MF C19 H25 N2 O4 . Br

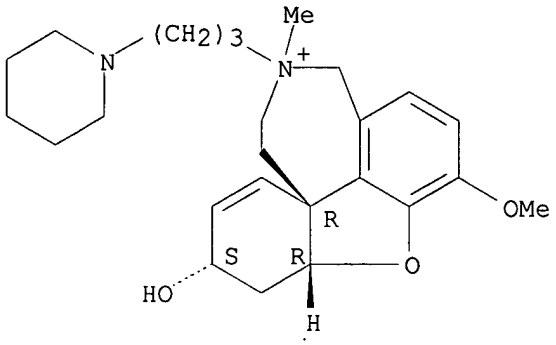
Absolute stereochemistry.



● Br^-

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-[3-(1-piperidinyl)propyl]-, chloride, (4aR,6S,8aR)-[partial]- (9CI)
MF C25 H37 N2 O3 . Cl

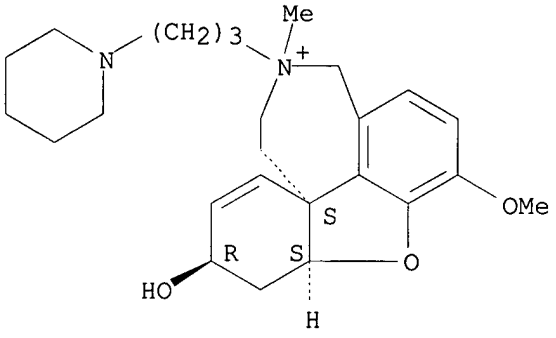
Absolute stereochemistry.



● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-[3-(1-piperidinyl)propyl]-, chloride, (4aS,6R,8aS)-[partial]- (9CI)
MF C25 H37 N2 O3 . Cl

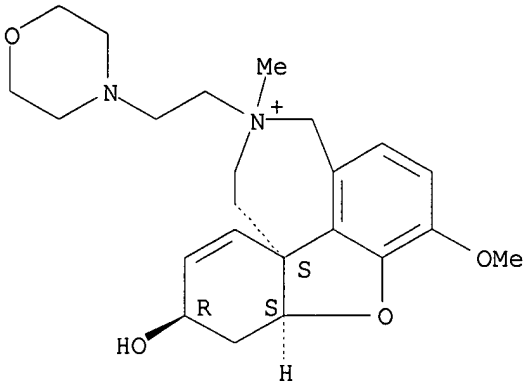
Absolute stereochemistry.



● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-[2-(4-morpholinyl)ethyl]-, chloride, (4aS,6R,8aS)-[partial]- (9CI)
MF C23 H33 N2 O4 . Cl

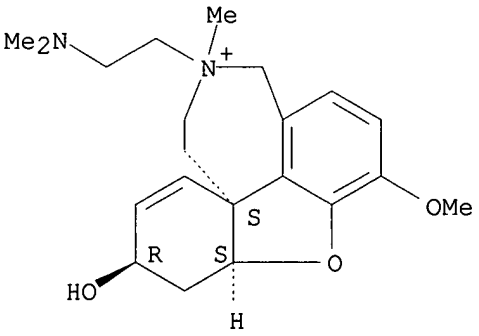
Absolute stereochemistry.



● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 11-[2-(dimethylamino)ethyl]-4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-, chloride, (4aS,6R,8aS)-[partial]- (9CI)
MF C21 H31 N2 O3 . Cl

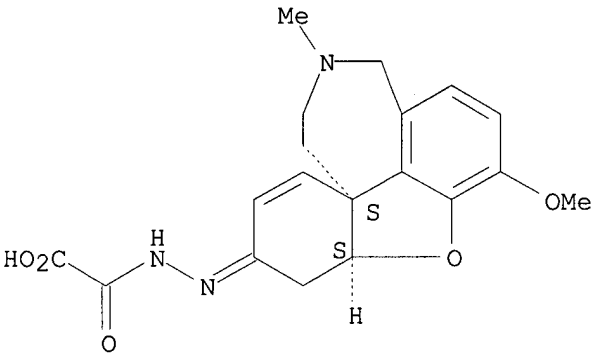
Absolute stereochemistry.



● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Ethanedioic acid, mono[(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)hydrazide], (4aR*,8aR*)-
(9CI)
MF C19 H21 N3 O5

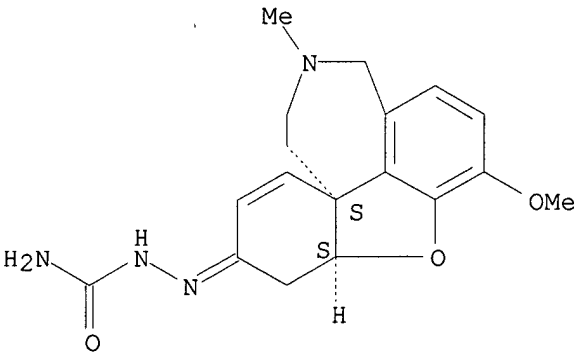
Relative stereochemistry.
Double bond geometry unknown.

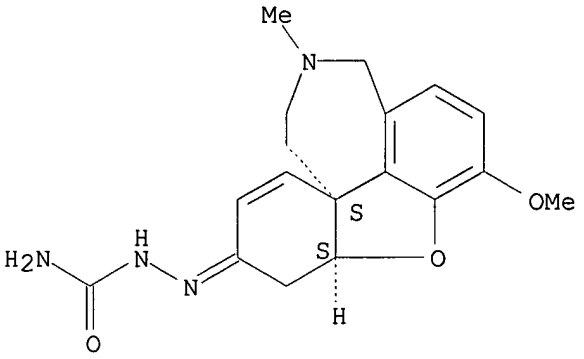


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazinecarboxamide,
2-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)-, (4aR*,8aR*)- (9CI)
MF C18 H22 N4 O3

Relative stereochemistry.
Double bond geometry unknown.

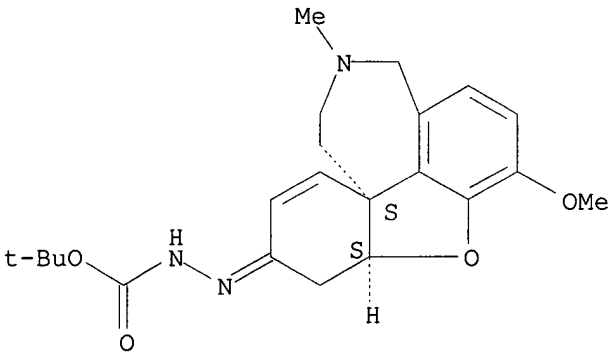




PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazinecarboxylic acid, (4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)-, 1,1-dimethylethyl ester,
(4aR*,8aR*)- (9CI)
MF C22 H29 N3 O4

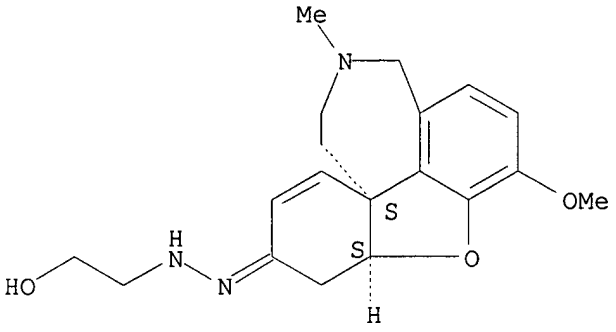
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (2-hydroxyethyl)hydrazone, [4aS-(4aR*,8aR*)]- (9CI)
MF C19 H25 N3 O3

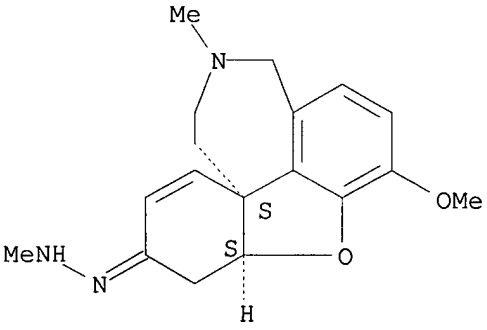
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, methylhydrazone, [4aS-(4aR*,8aR*)]- (9CI)
MF C18 H23 N3 O2

Absolute stereochemistry.
Double bond geometry unknown.

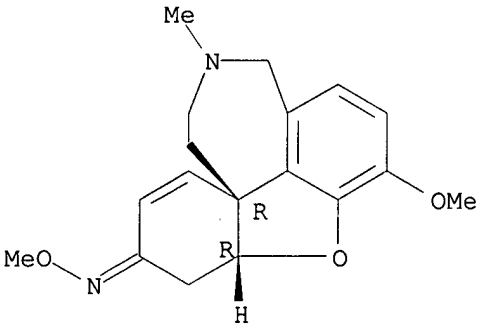


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-

methoxy-11-methyl-, O-methyloxime, [4aR-(4aR*,8aR*)]- (9CI)
MF C18 H22 N2 O3

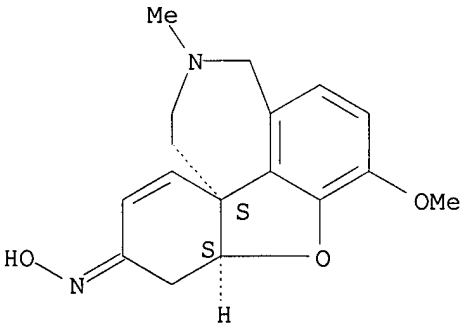
Absolute stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, oxime, [4aS-(4aR*,8aR*)]- (9CI)
MF C17 H20 N2 O3

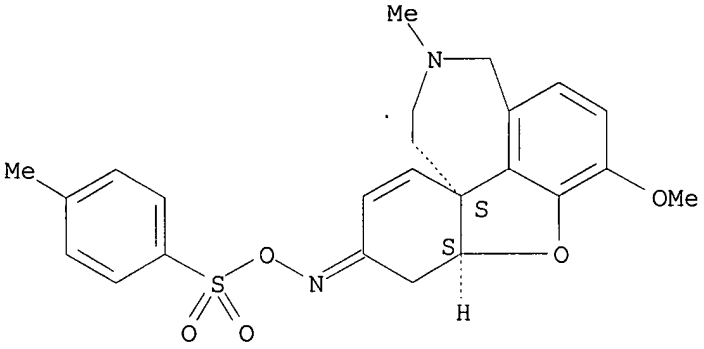
Absolute stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, O-[(4-methylphenyl)sulfonyl]oxime, (4aR*,8aR*)- (9CI)
MF C24 H26 N2 O5 S

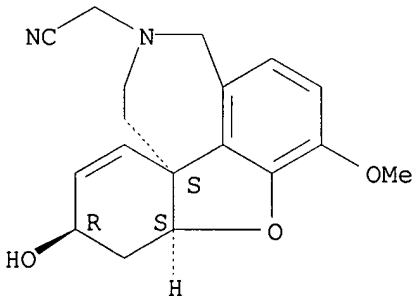
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetonitrile, 4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, (4aR,6S,8aR)-rel- (9CI)
MF C18 H20 N2 O3

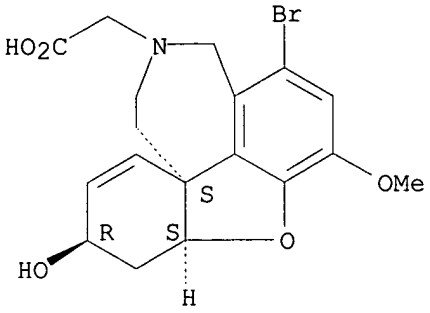
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetic acid, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H20 Br N O5

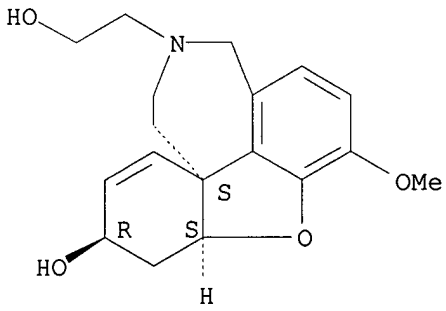
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-ethanol, 4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, (4aR,6S,8aR)-rel- (9CI)
MF C18 H23 N O4

Relative stereochemistry.

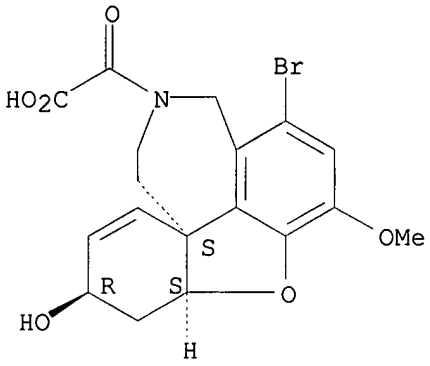


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetic acid, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-.alpha.-oxo-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H18 Br N O6

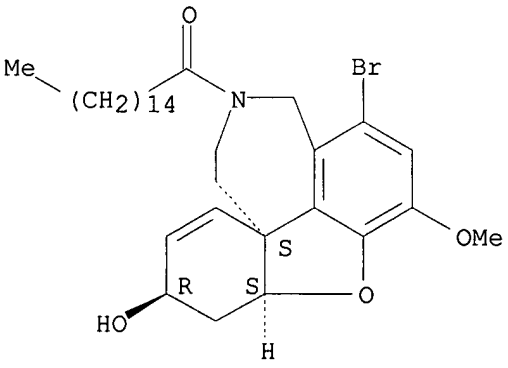
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-(1-oxohexadecyl)-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C32 H48 Br N O4

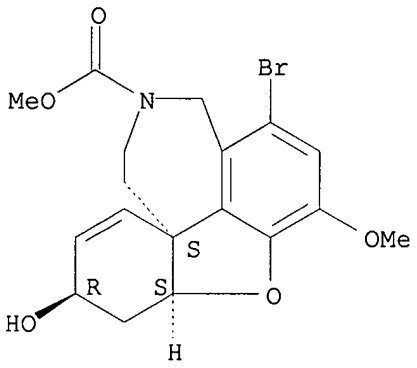
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-carboxylic acid, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, methyl ester, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H20 Br N O5

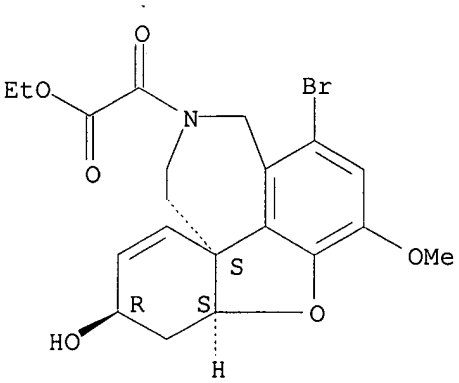
Relative stereochemistry.

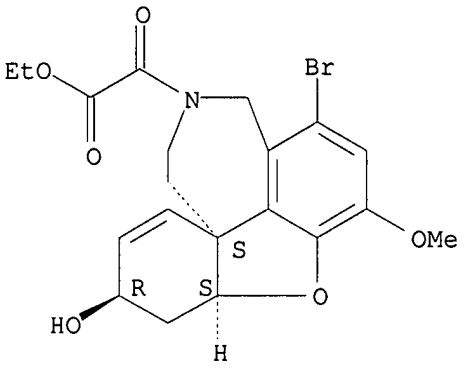


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetic acid, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-.alpha.-oxo-, ethyl ester, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C20 H22 Br N O6

Relative stereochemistry.

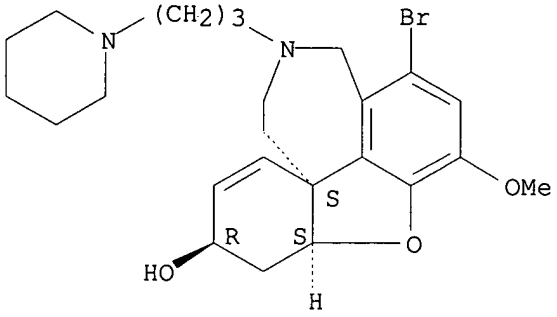




PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-[3-(1-piperidinyl)propyl]-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C24 H33 Br N2 O3

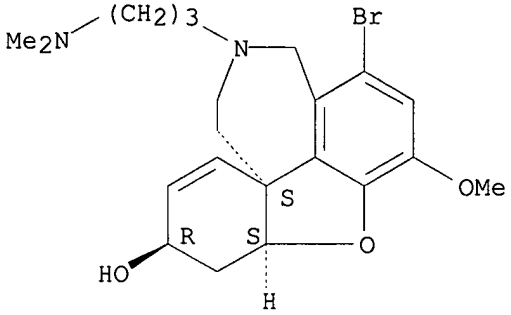
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-11-[3-(dimethylamino)propyl]-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C21 H29 Br N2 O3

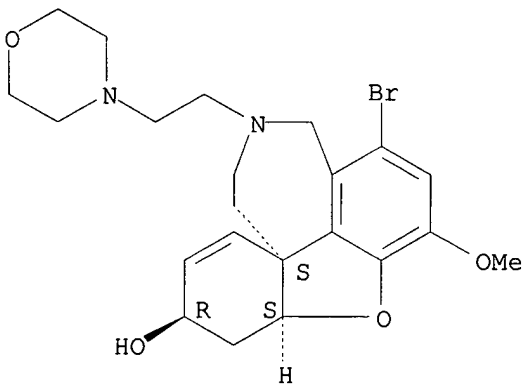
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-[2-(4-morpholinyl)ethyl]-, (4a.alpha.,6.beta.,8aR*)-
(9CI)
MF C22 H29 Br N2 O4

Relative stereochemistry.

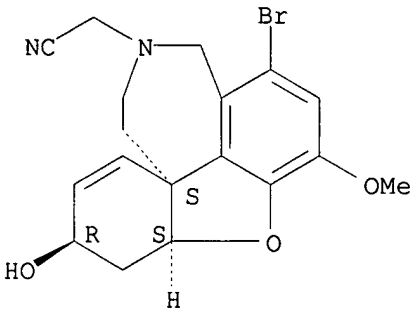


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetonitrile, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)

MF C18 H19 Br N2 O3

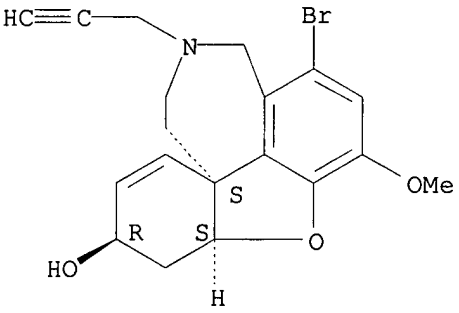
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-(2-propynyl)-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C19 H20 Br N O3

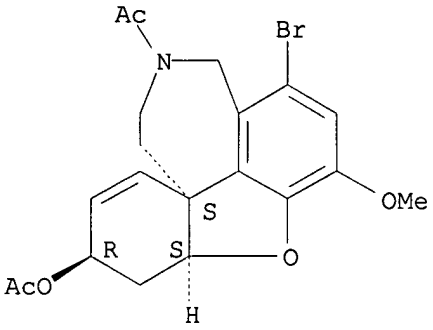
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 11-acetyl-1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-, acetate (ester), (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C20 H22 Br N O5

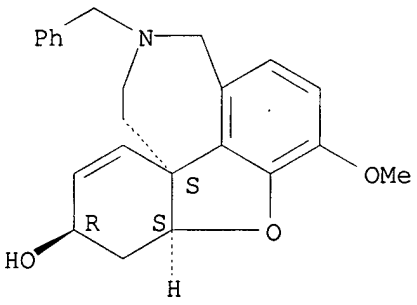
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-(phenylmethyl)-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C23 H25 N O3

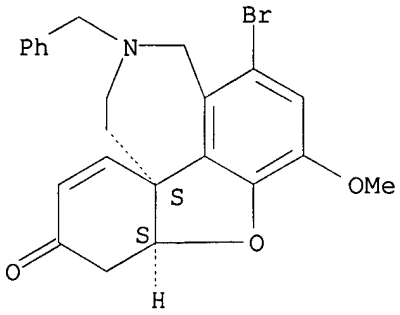
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-(phenylmethyl)-, (4aR*,8aR*)- (9CI)
MF C23 H22 Br N O3

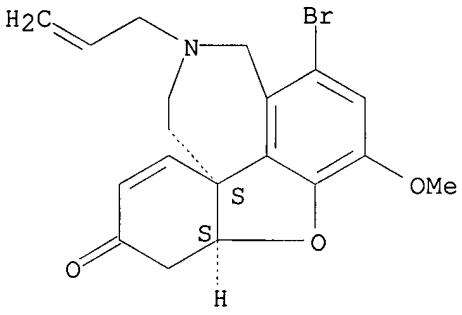
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-(2-propenyl)-, (4aR*,8aR*)- (9CI)
MF C19 H20 Br N O3

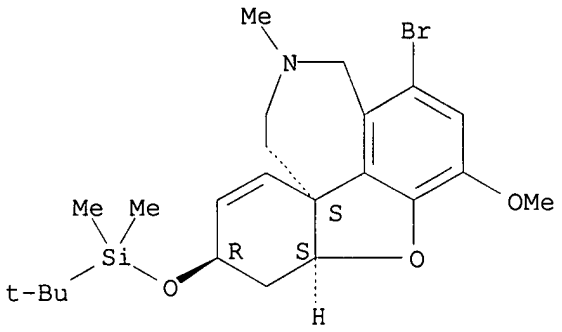
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine, 1-bromo-6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, [4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C23 H34 Br N O3 Si

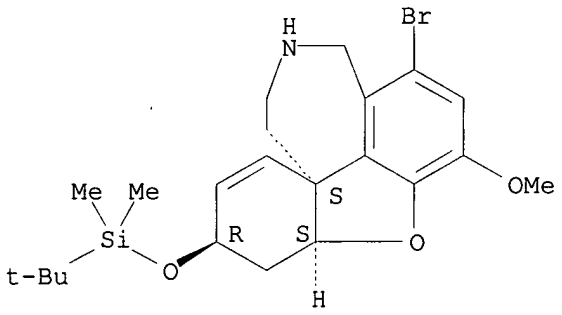
Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine, 1-bromo-6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C22 H32 Br N O3 Si

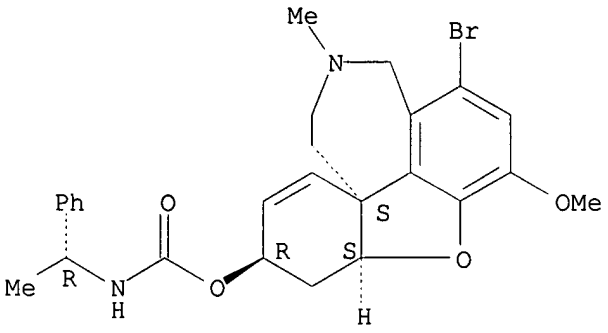
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, (1-phenylethyl)-, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester, [4a.alpha.,6.beta.(S*),8aR*]- (9CI)
MF C26 H29 Br N2 O4

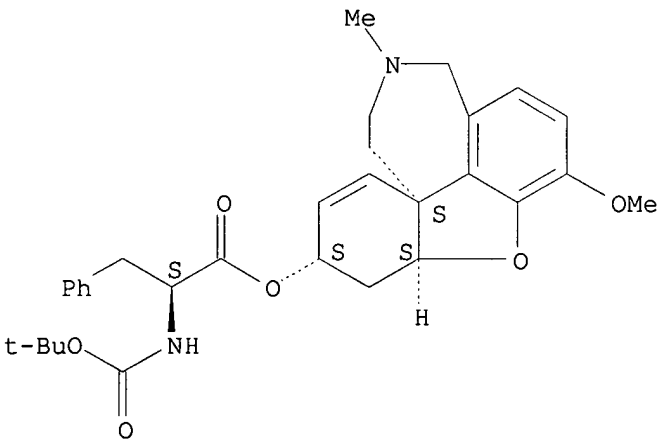
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Phenylalanine, N-[(1,1-dimethylethoxy)carbonyl]-, (4aS,6S,8aS)-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester (9CI)
MF C31 H38 N2 O6

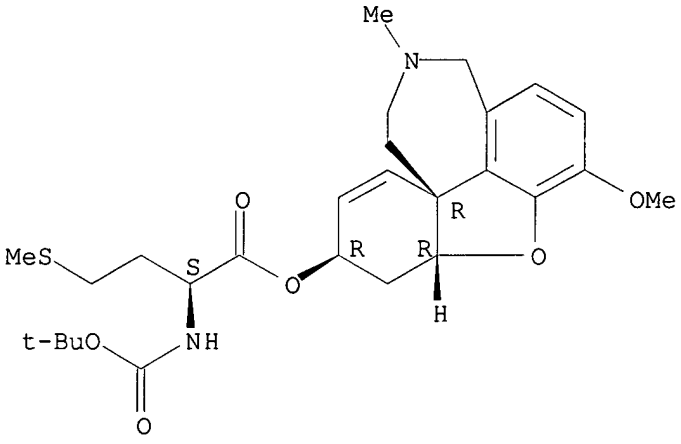
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Methionine, N-[(1,1-dimethylethoxy)carbonyl]-, (4aR,6R,8aR)-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester (9CI)
MF C27 H38 N2 O6 S

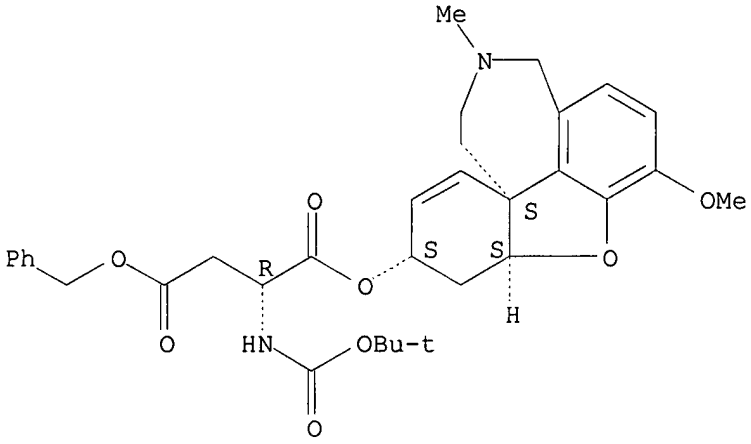
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN D-Aspartic acid, N-[(1,1-dimethylethoxy)carbonyl]-, 1-[(4aS,6S,8aS)-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl] 4-(phenylmethyl) ester (9CI)
MF C33 H40 N2 O8

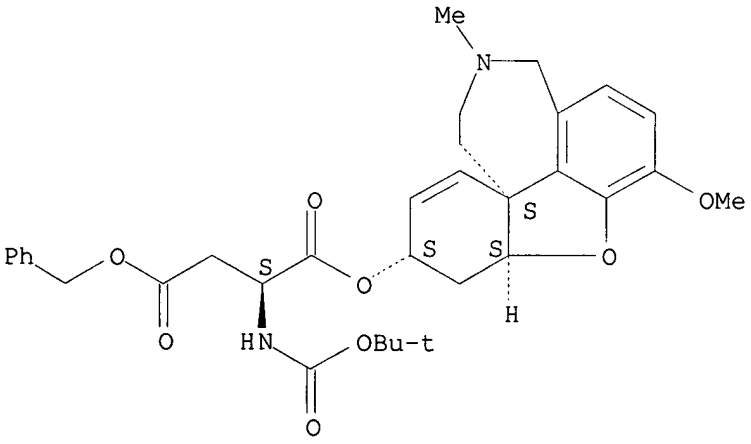
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Aspartic acid, N-[(1,1-dimethylethoxy)carbonyl]-, 1-[(4aS,6S,8aS)-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl] 4-(phenylmethyl) ester (9CI)
MF C33 H40 N2 O8

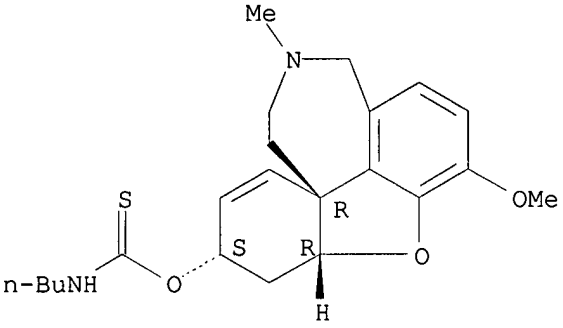
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamothioic acid, butyl-, O-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl) ester, [4aR-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C22 H30 N2 O3 S

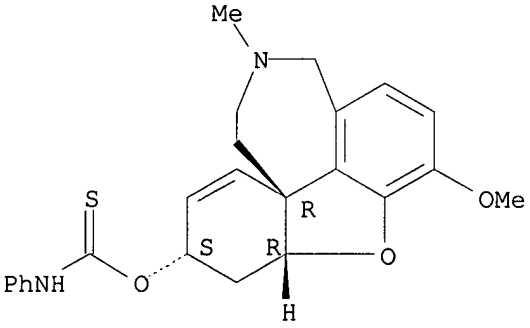
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamothioic acid, phenyl-, O-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl) ester, [4aR-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C24 H26 N2 O3 S

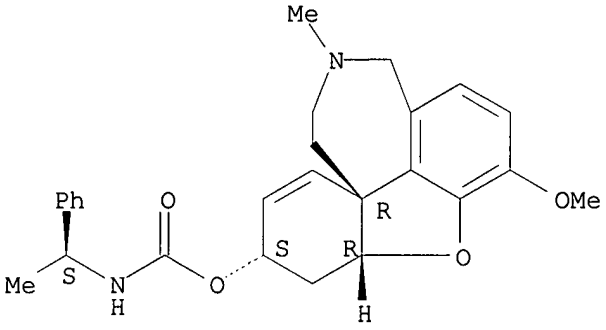
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, (1-phenylethyl)-, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester, [4aR-[4a.alpha.,6.beta.(S*),8aR*]]- (9CI)
MF C26 H30 N2 O4

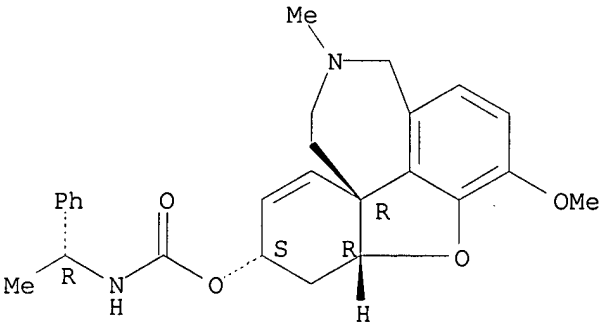
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, (1-phenylethyl)-, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester, [4aR-[4a.alpha.,6.beta.(R*),8aR*]]- (9CI)
MF C26 H30 N2 O4

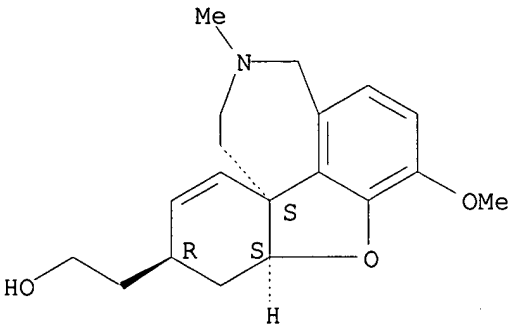
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-6-ethanol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C19 H25 N O3

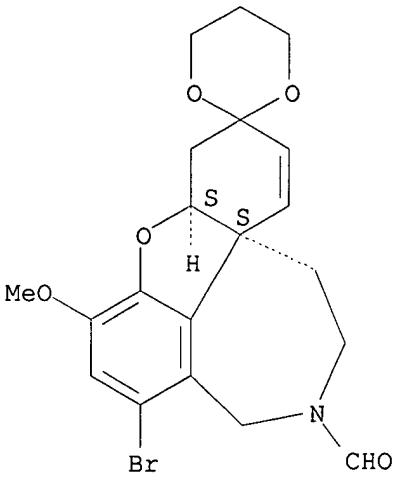
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxane]-11(12H)-carboxaldehyde, 1-bromo-4a,5,9,10-tetrahydro-3-methoxy-, (4aR*,8aR*)- (9CI)
MF C20 H22 Br N O5

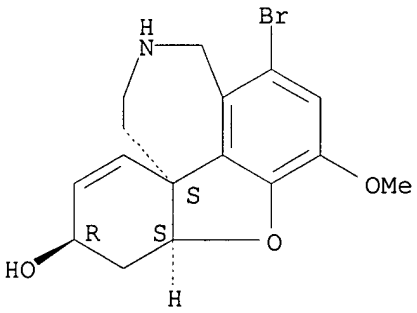
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4aS,6R,8aS)- (9CI)
MF C16 H18 Br N O3

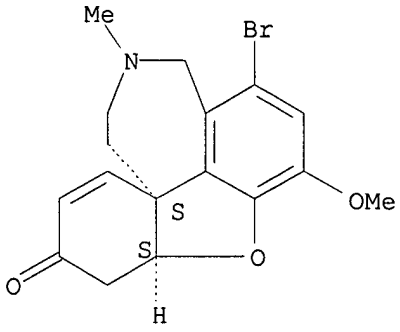
Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR*,8aR*)- (9CI)
MF C17 H18 Br N O3

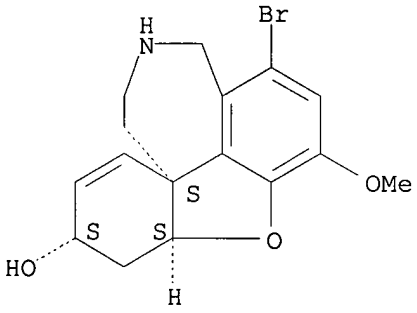
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4aR,6R,8aR)-rel- (9CI)
MF C16 H18 Br N O3

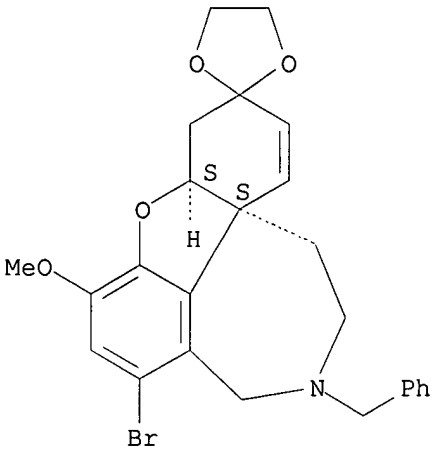
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxolane], 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-(phenylmethyl)-, (4aR*,8aR*)- (9CI)
MF C25 H26 Br N O4

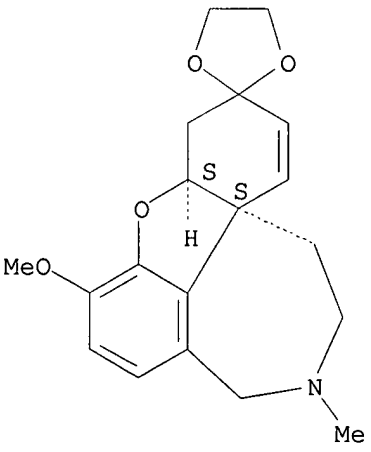
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxolane], 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR*,8aR*)- (9CI)
MF C19 H23 N O4

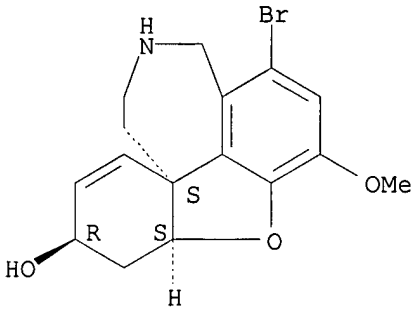
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4aR,6S,8aR)-rel- (9CI)
MF C16 H18 Br N O3

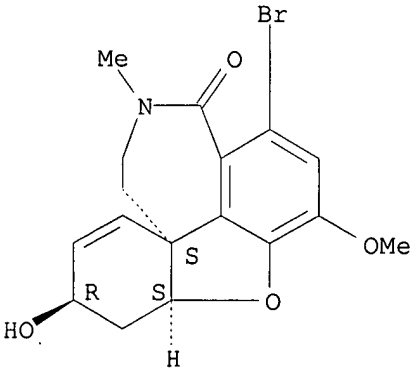
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

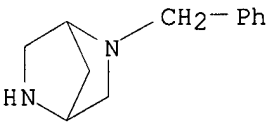
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-12(9H)-one, 1-bromo-4a,5,10,11-tetrahydro-6-hydroxy-3-methoxy-11-methyl-, (4aR,6S,8aR)-rel- (9CI)
MF C17 H18 Br N O4

Relative stereochemistry.



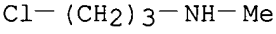
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(phenylmethyl)-, dihydrobromide (9CI)
MF C12 H16 N2 . 2 Br H



●2 HBr

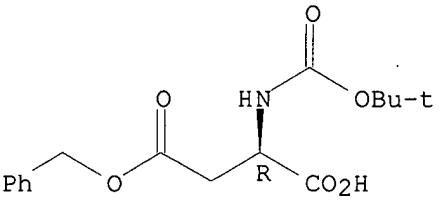
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 1-Propanamine, 3-chloro-N-methyl- (9CI)
MF C4 H10 Cl N
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

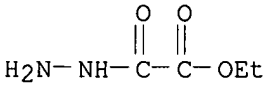
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN D-Aspartic acid, N-[(1,1-dimethylethoxy)carbonyl]-, 4-(phenylmethyl) ester
(9CI)
MF C16 H21 N O6

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

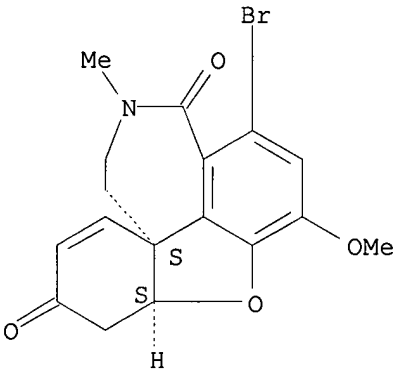
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Ethanedioic acid, monoethyl ester, hydrazide (9CI)
MF C4 H8 N2 O3
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Galanthamine, 8-bromo-3-deoxy-3,9-dioxo- (9CI)
MF C17 H16 Br N O4

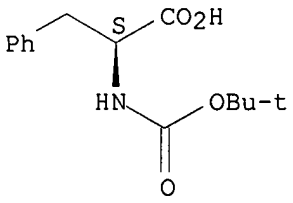
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Phenylalanine, N-[(1,1-dimethylethoxy)carbonyl]- (9CI)
MF C14 H19 N O4
CI COM

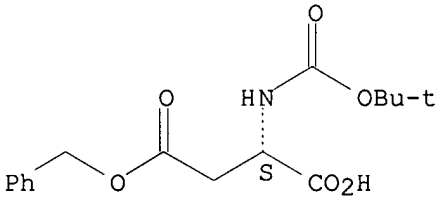
Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

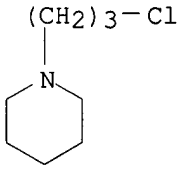
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Aspartic acid, N-[(1,1-dimethylethoxy)carbonyl]-, 4-(phenylmethyl) ester
(9CI)
MF C16 H21 N O6
CI COM

Absolute stereochemistry.



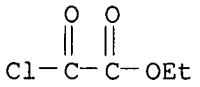
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Piperidine, 1-(3-chloropropyl)-, hydrochloride (6CI, 7CI, 8CI, 9CI)
MF C8 H16 Cl N . Cl H



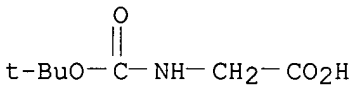
● HCl

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Acetic acid, chlorooxo-, ethyl ester (9CI)
MF C4 H5 Cl O3



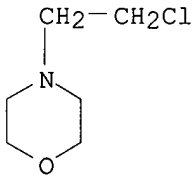
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Glycine, N-[(1,1-dimethylethoxy)carbonyl]- (9CI)
MF C7 H13 N O4
CI COM



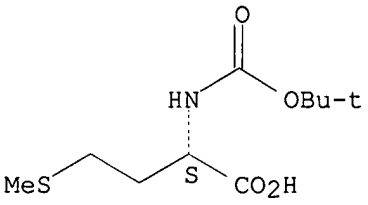
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Morpholine, 4-(2-chloroethyl)-, hydrochloride (6CI, 7CI, 8CI, 9CI)
MF C6 H12 Cl N O . Cl H



L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Methionine, N-[(1,1-dimethylethoxy)carbonyl]- (9CI)
MF C10 H19 N O4 S
CI COM

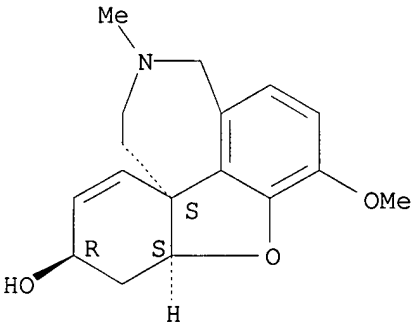
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

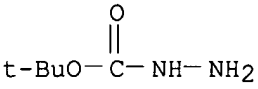
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, hydrobromide, (4aS,6R,8aS)- (9CI)
MF C17 H21 N O3 . Br H
CI COM

Absolute stereochemistry. Rotation (-).



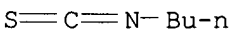
● HBr

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazinecarboxylic acid, 1,1-dimethylethyl ester (9CI)
MF C5 H12 N2 O2
CI COM



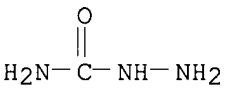
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Butane, 1-isothiocyanato- (9CI)
MF C5 H9 N S
CI COM



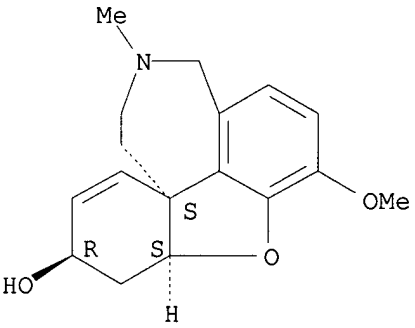
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazinecarboxamide, monohydrochloride (9CI)
MF C H5 N3 O . Cl H
CI COM



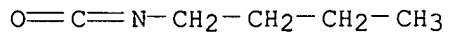
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aS,6R,8aS)- (9CI)
MF C17 H21 N O3
CI COM

Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Butane, 1-isocyanato- (9CI)
MF C5 H9 N O
CI COM

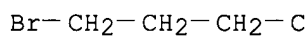


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Pentane, 1-bromo- (6CI, 8CI, 9CI)
MF C5 H11 Br
CI COM

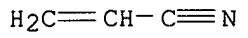
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Propane, 1-bromo-3-chloro- (6CI, 7CI, 8CI, 9CI)
MF C3 H6 Br Cl
CI COM



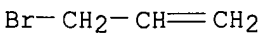
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2-Propenenitrile (9CI)
MF C3 H3 N
CI COM



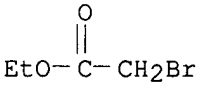
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 1-Propene, 3-bromo- (9CI)
MF C3 H5 Br
CI COM



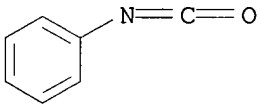
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Acetic acid, bromo-, ethyl ester (6CI, 8CI, 9CI)
MF C4 H7 Br O2
CI COM



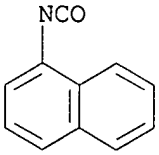
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzene, isocyanato- (9CI)
MF C7 H5 N O
CI COM



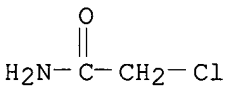
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Naphthalene, 1-isocyanato- (9CI)
MF C11 H7 N O
CI COM



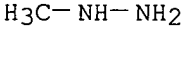
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Acetamide, 2-chloro- (6CI, 7CI, 8CI, 9CI)
MF C2 H4 Cl N O
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

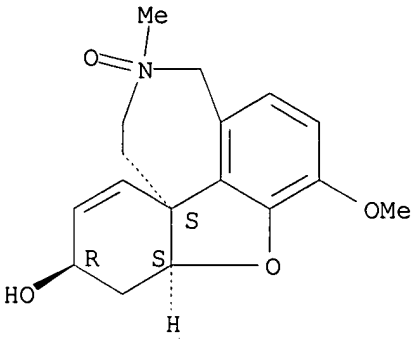
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazine, methyl- (6CI, 8CI, 9CI)
MF C H6 N2
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

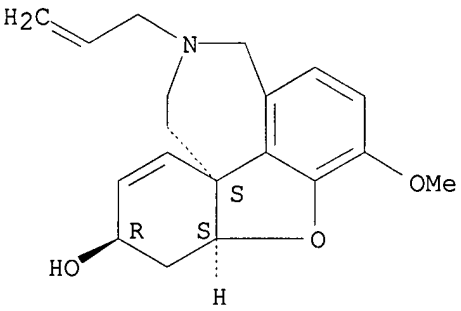
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, 11-oxide, [4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C17 H21 N O4

Absolute stereochemistry.



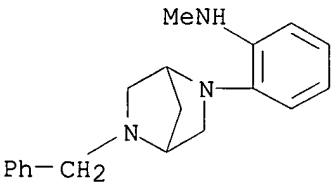
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-(2-propenyl)-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C19 H23 N O3

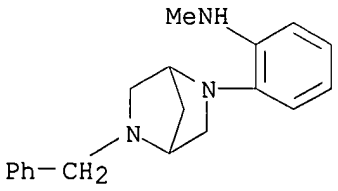
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

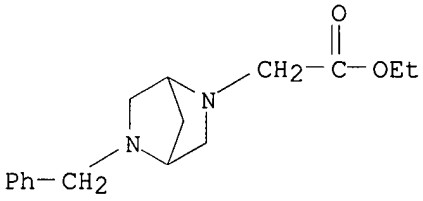
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenamine, N-methyl-2-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]- (9CI)
MF C19 H23 N3





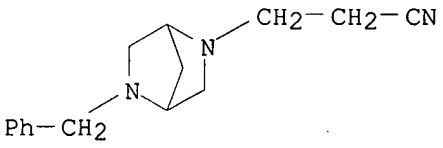
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane-2-acetic acid, 5-(phenylmethyl)-, ethyl ester (9CI)
MF C16 H22 N2 O2



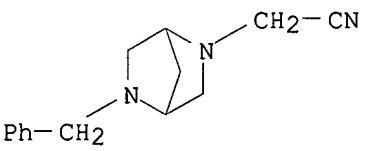
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane-2-propanenitrile, 5-(phenylmethyl)- (9CI)
MF C15 H19 N3



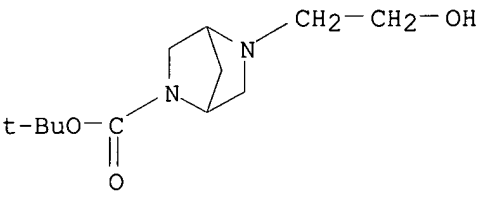
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane-2-acetonitrile, 5-(phenylmethyl)- (9CI)
MF C14 H17 N3



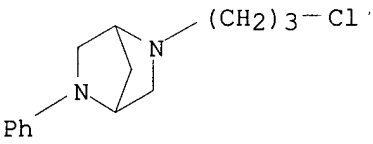
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane-2-carboxylic acid, 5-(2-hydroxyethyl)-, 1,1-dimethylethyl ester (9CI)
MF C12 H22 N2 O3



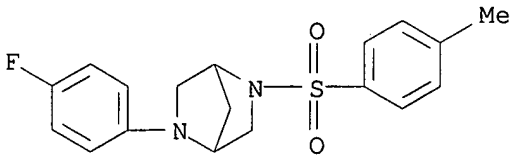
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(3-chloropropyl)-5-phenyl- (9CI)
MF C14 H19 Cl N2



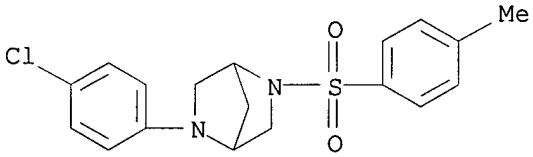
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-fluorophenyl)-5-[(4-methylphenyl)sulfonyl]- (9CI)
MF C18 H19 F N2 O2 S



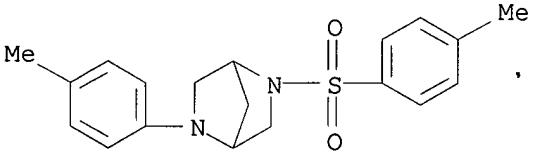
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-chlorophenyl)-5-[(4-methylphenyl)sulfonyl]- (9CI)
MF C18 H19 Cl N2 O2 S



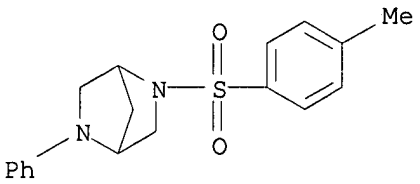
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-methylphenyl)-5-[(4-methylphenyl)sulfonyl]- (9CI)
MF C19 H22 N2 O2 S



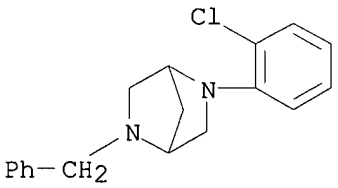
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-[(4-methylphenyl)sulfonyl]-5-phenyl- (9CI)
MF C18 H20 N2 O2 S



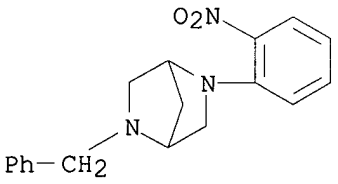
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(2-chlorophenyl)-5-(phenylmethyl)-(9CI)
MF C18 H19 Cl N2



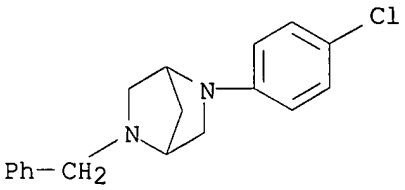
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(2-nitrophenyl)-5-(phenylmethyl)- (9CI)
MF C18 H19 N3 O2



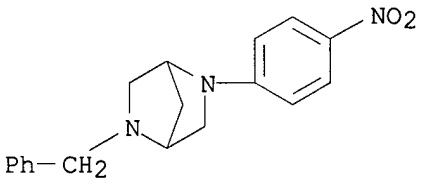
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-chlorophenyl)-5-(phenylmethyl)-(9CI)
MF C18 H19 Cl N2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

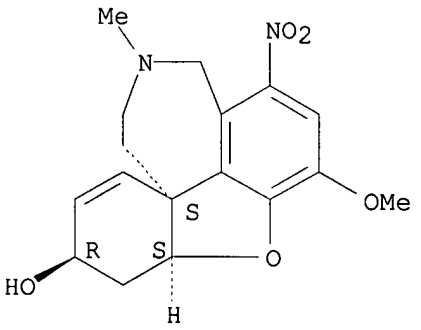
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-nitrophenyl)-5-(phenylmethyl)- (9CI)
MF C18 H19 N3 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-1-nitro-, [4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C17 H20 N2 O5

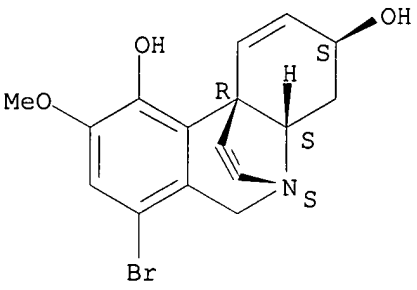
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

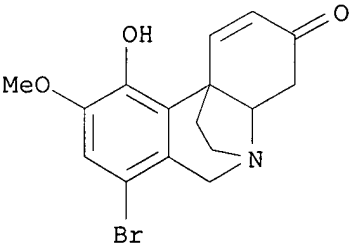
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 3H,6H-5,10b-Ethynophenanthridine-3,10-diol, 7-bromo-4,4a-dihydro-9-methoxy-
, (3.alpha.,4a.alpha.,5.alpha.,10b.alpha.)- (9CI)
MF C16 H14 Br N O3

Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 3H,6H-5,10b-Ethanophenanthridin-3-one, 7-bromo-4,4a-dihydro-10-hydroxy-9-methoxy- (9CI)
MF C16 H16 Br N O3

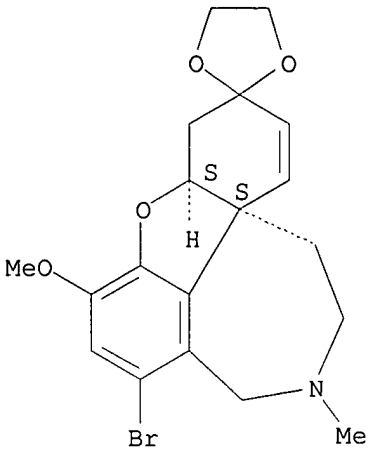


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxolane], 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR*,8aR*)-
(9CI)

MF C19 H22 Br N O4

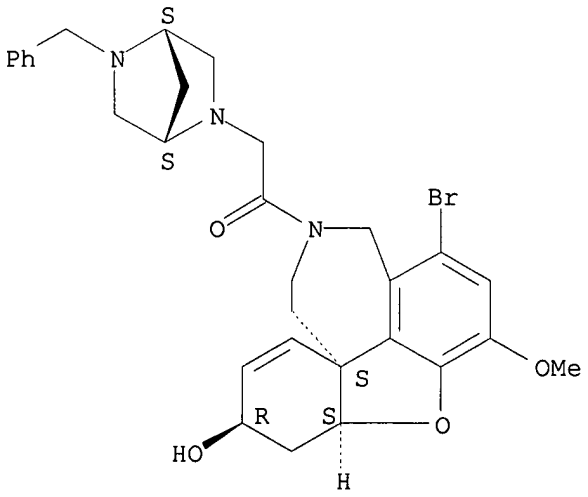
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-[[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]acetyl]-, [4a.alpha.,6.beta.,8aR*,11(1R*,4R*)]- (9CI)
MF C30 H34 Br N3 O4

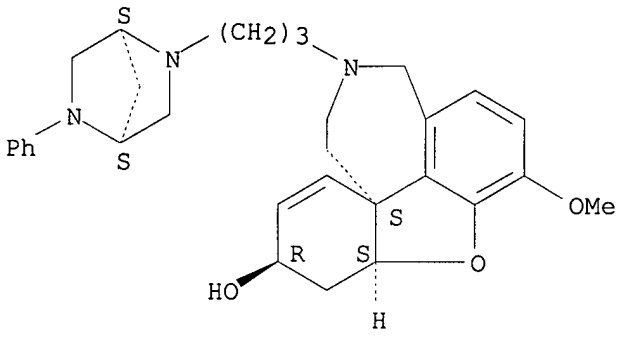
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-[3-(5-phenyl-2,5-diazabicyclo[2.2.1]hept-2-yl)propyl]-, [4a.alpha.,6.beta.,8aR*,11(1R*,4R*)]- (9CI)
MF C30 H37 N3 O3

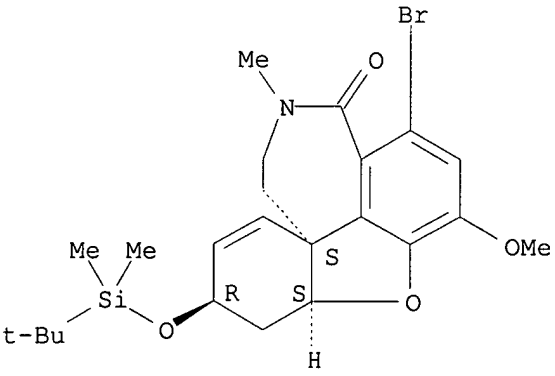
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-12(9H)-one, 1-bromo-6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-4a,5,10,11-tetrahydro-3-methoxy-11-methyl-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C23 H32 Br N O4 Si

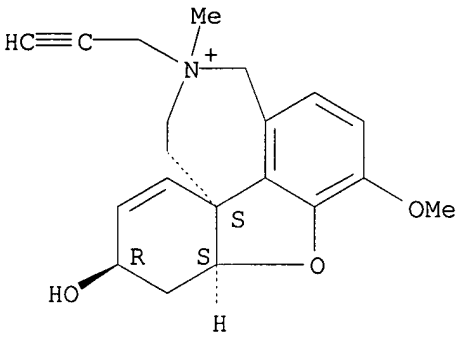
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-(2-propynyl)-, bromide, (4aS,6R,8aS)-[partial]- (9CI)
MF C20 H24 N O3 . Br

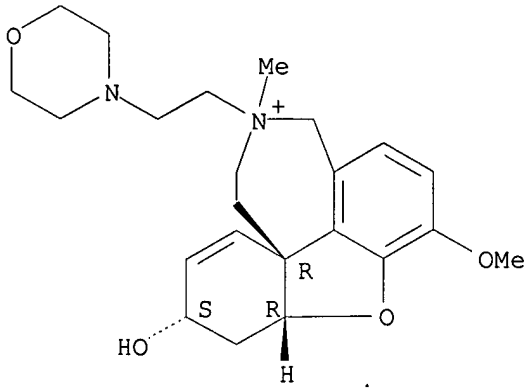
Absolute stereochemistry.



● Br^-

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-[2-(4-morpholinyl)ethyl]-, chloride, (4aR,6S,8aR)-[partial]- (9CI)
MF C23 H33 N2 O4 . Cl

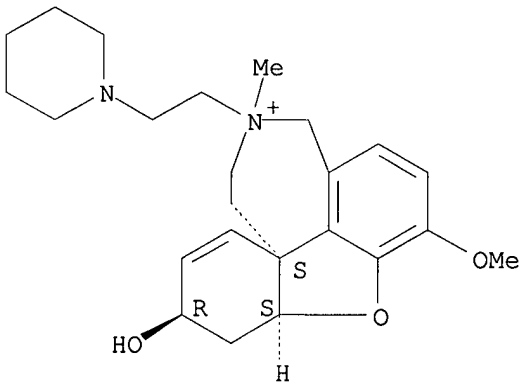
Absolute stereochemistry.



● Cl^-

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-[2-(1-piperidinyl)ethyl]-, chloride, (4aS,6R,8aS)-[partial]- (9CI)
MF C24 H35 N2 O3 . Cl

Absolute stereochemistry.

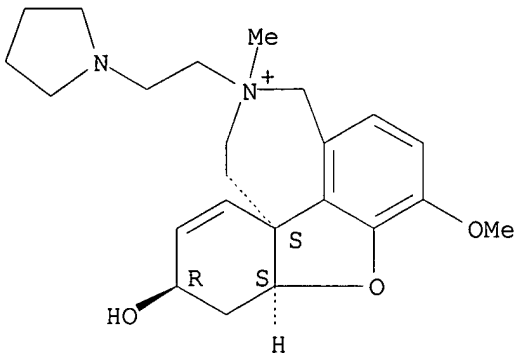


● Cl^-

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-

hydroxy-3-methoxy-11-methyl-11-[2-(1-pyrrolidinyl)ethyl]-, chloride, (4aS,6R,8aS)-[partial]- (9CI)
MF C23 H33 N2 O3 . Cl

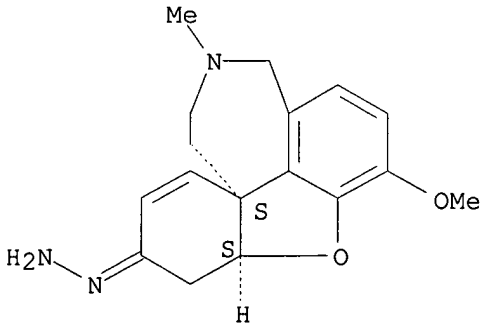
Absolute stereochemistry.



● Cl^-

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, hydrazone, [4aS-(4aR*,8aR*)]- (9CI)
MF C17 H21 N3 O2

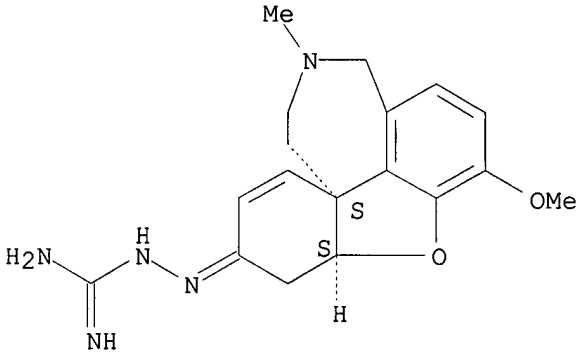
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazinecarboximidamide,
2-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-
6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)-, (4aR*,8aR*)- (9CI)
MF C18 H23 N5 O2

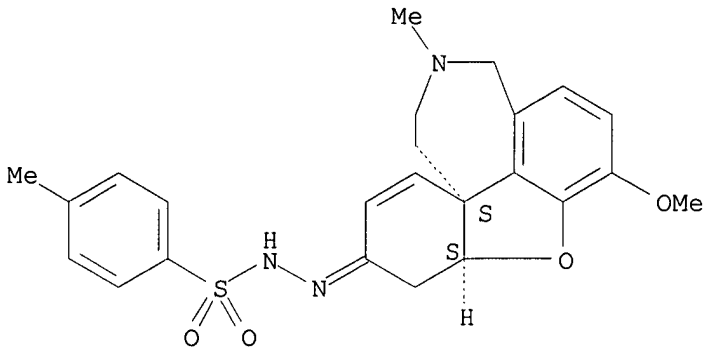
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenesulfonic acid, 4-methyl-, (4a,5,9,10,11,12-hexahydro-3-methoxy-11-
methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)hydrazide,
(4aR*,8aR*)- (9CI)
MF C24 H27 N3 O4 S

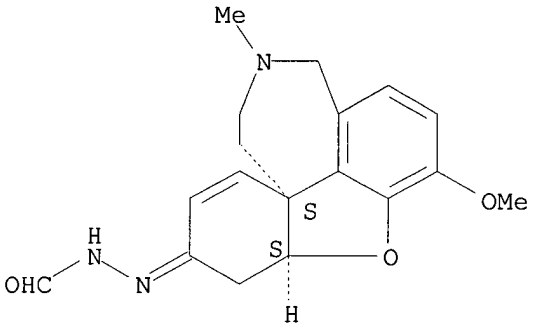
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazinecarboxaldehyde,
(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-
benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)-, (4aR*,8aR*)- (9CI)
MF C18 H21 N3 O3

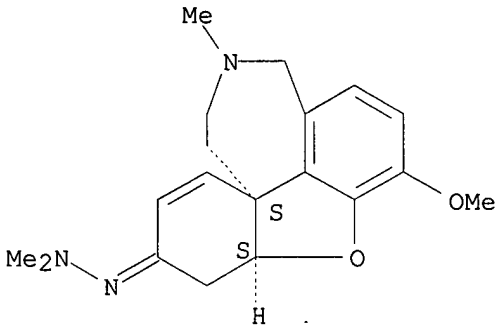
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-
methoxy-11-methyl-, dimethylhydrazone, [4aS-(4aR*,8aR*)]- (9CI)
MF C19 H25 N3 O2

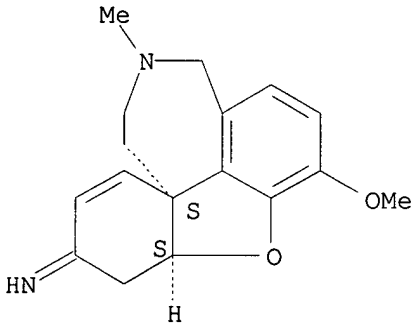
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-imine, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, [4aS-(4aR*,8aR*)]- (9CI)
MF C17 H20 N2 O2

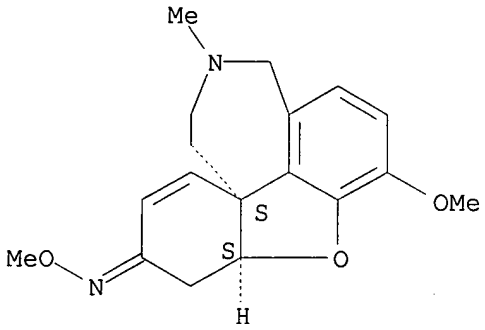
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, O-methyloxime, [4aS-(4aR*,8aR*)]- (9CI)
MF C18 H22 N2 O3

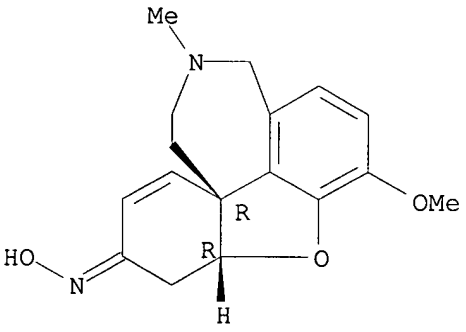
Absolute stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, oxime, [4aR-(4aR*,8aR*)]- (9CI)
MF C17 H20 N2 O3

Absolute stereochemistry.
Double bond geometry unknown.

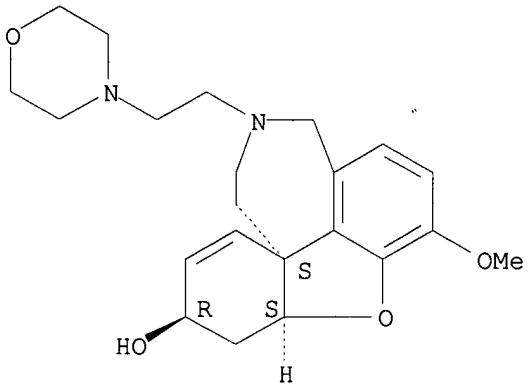


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-[2-(4-morpholinyl)ethyl]-, (4a.alpha.,6.beta.,8aR*)- (9CI)

MF C22 H30 N2 O4

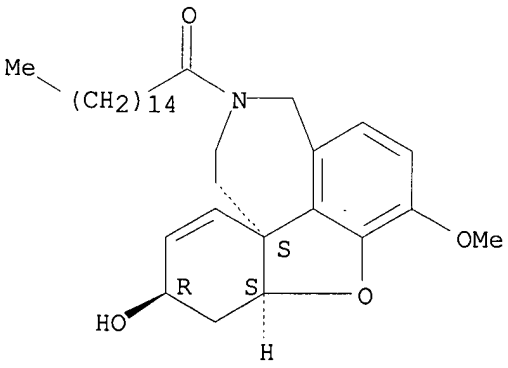
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-(1-oxohexadecyl)-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C32 H49 N O4

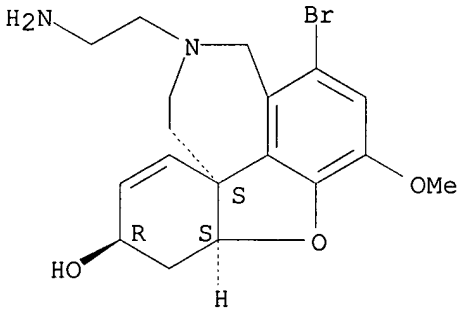
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 11-(2-aminoethyl)-1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H23 Br N2 O3

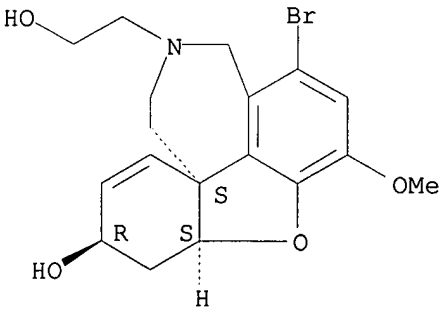
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-ethanol, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H22 Br N O4

Relative stereochemistry.

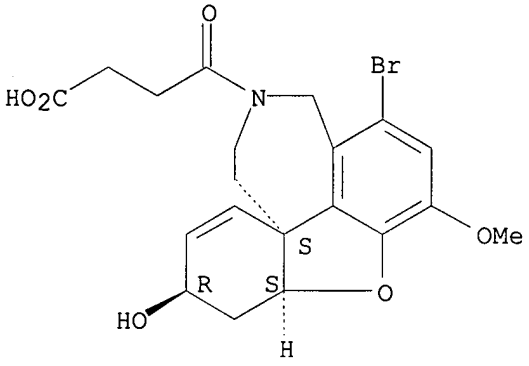


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-butanoic acid, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-.gamma.-oxo-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C20 H22 Br N O6

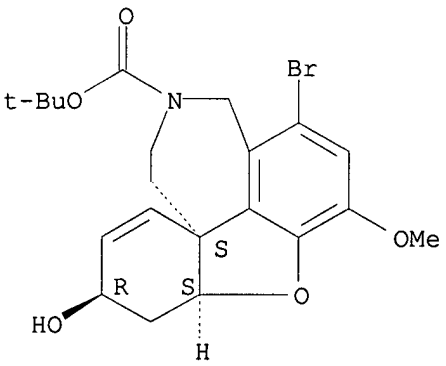
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-carboxylic acid, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, 1,1-dimethylethyl ester, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C21 H26 Br N O5

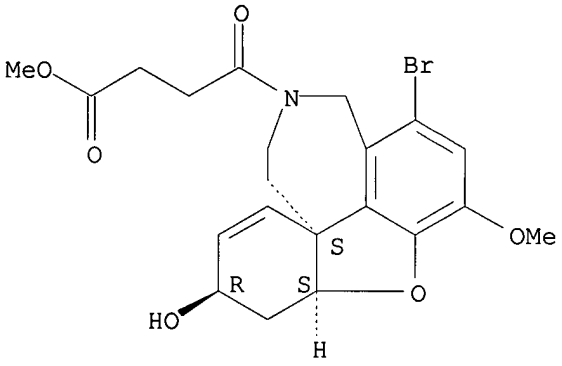
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-butanoic acid, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-.gamma.-oxo-, methyl ester, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C21 H24 Br N O6

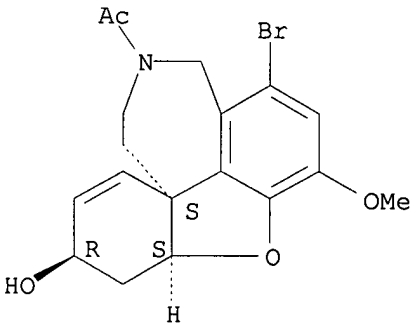
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 11-acetyl-1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H20 Br N O4

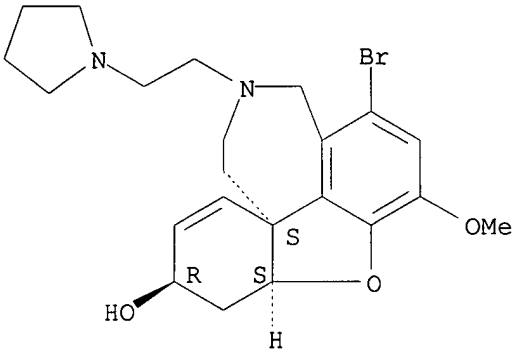
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-[2-(1-pyrrolidinyl)ethyl]-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C22 H29 Br N2 O3

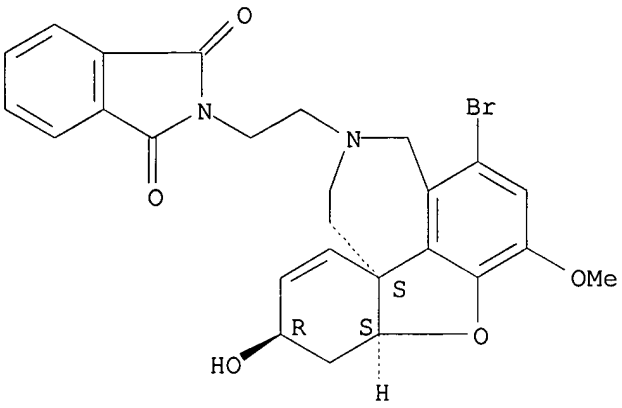
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 1H-Isoindole-1,3(2H)-dione, 2-[2-(1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-6H-benzofuro[3a,3,2-ef][2]benzazepin-11(12H)-yl)ethyl]-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C26 H25 Br N2 O5

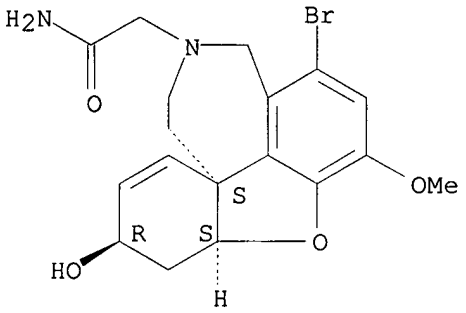
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetamide, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H21 Br N2 O4

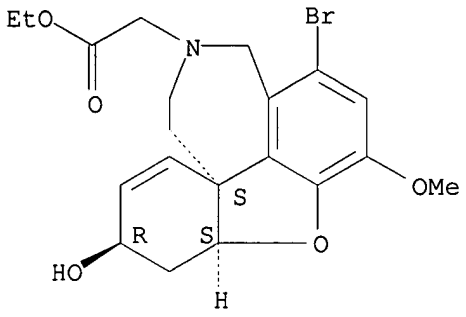
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetic acid, 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, ethyl ester, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C20 H24 Br N O5

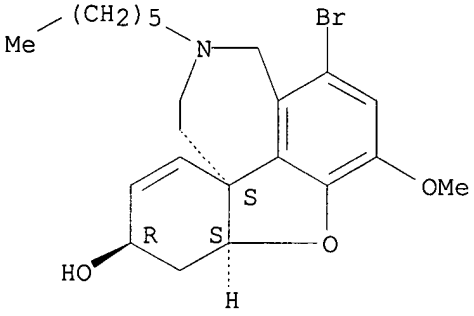
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-11-hexyl-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C22 H30 Br N O3

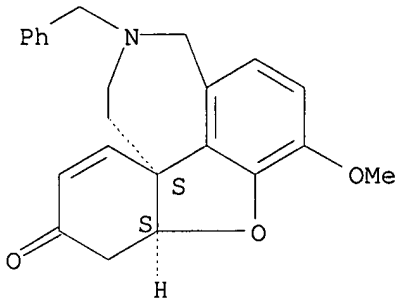
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-(phenylmethyl)-, (4aR*,8aR*)- (9CI)
MF C23 H23 N O3

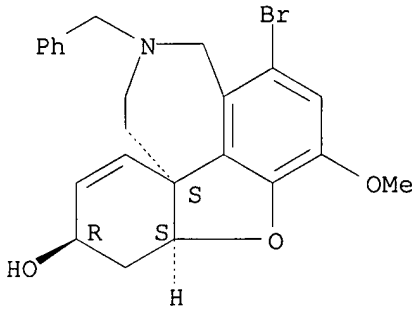
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-(phenylmethyl)-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C23 H24 Br N O3

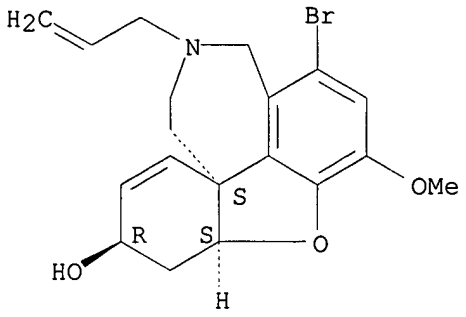
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-(2-propenyl)-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C19 H22 Br N O3

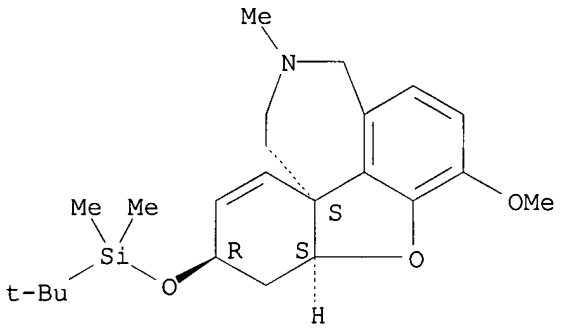
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine, 6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, [4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C23 H35 N O3 Si

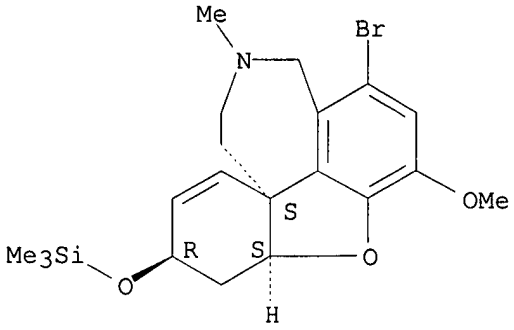
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6-[(trimethylsilyl)oxy]-, (4a.alpha.,6.beta.,8aR*)-(9CI)
MF C20 H28 Br N O3 Si

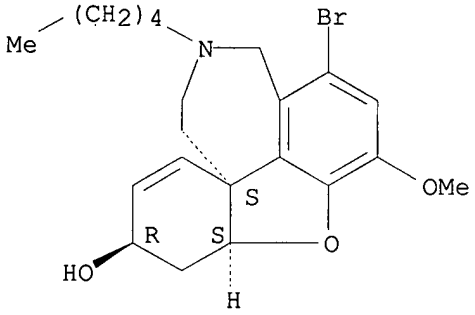
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-pentyl-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C21 H28 Br N O3

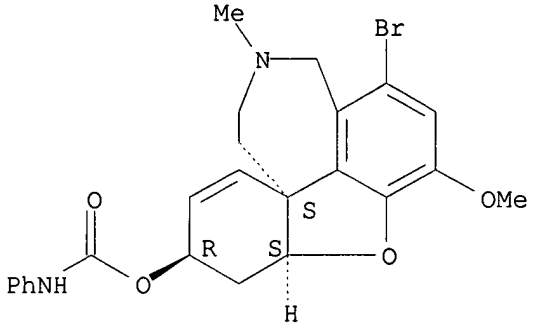
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, phenylcarbamate (ester), (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C24 H25 Br N2 O4

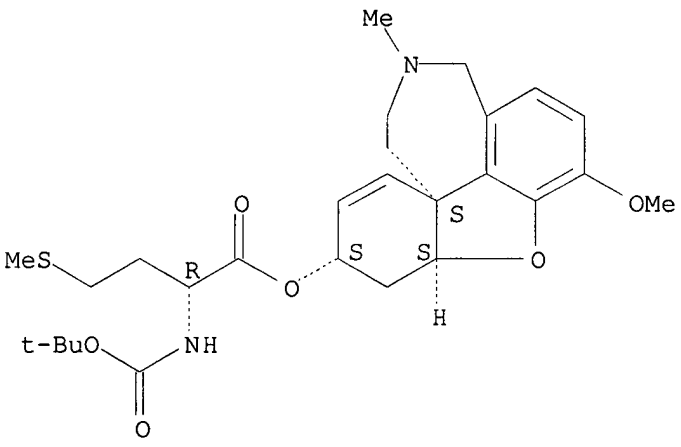
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN D-Methionine, N-[(1,1-dimethylethoxy)carbonyl]-, (4aS,6S,8aS)-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester (9CI)
MF C27 H38 N2 O6 S

Absolute stereochemistry.

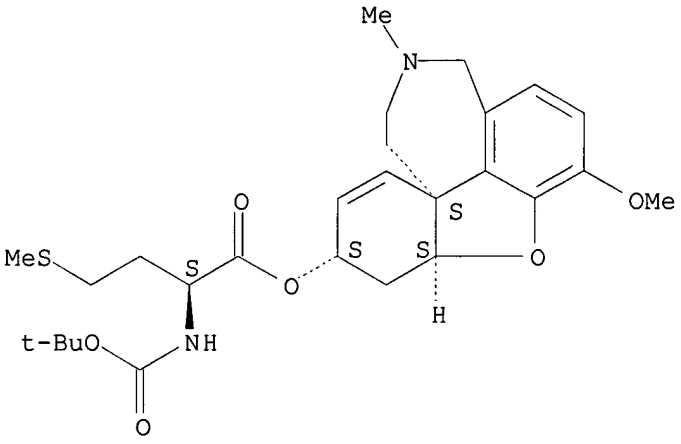


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Methionine, N-[(1,1-dimethylethoxy)carbonyl]-, (4aS,6S,8aS)-

4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester (9CI)
MF C27 H38 N2 O6 S

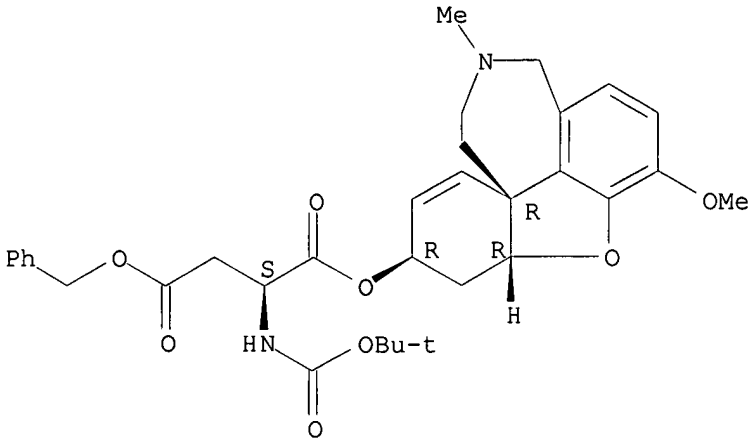
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Aspartic acid, N-[(1,1-dimethylethoxy)carbonyl]-, 1-[(4aR,6R,8aR)-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl] 4-(phenylmethyl) ester (9CI)
MF C33 H40 N2 O8

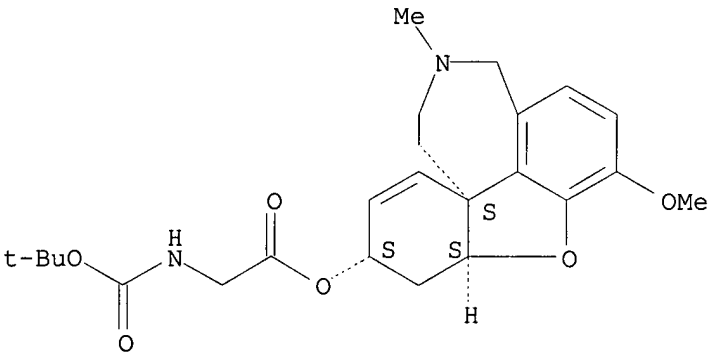
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Glycine, N-[(1,1-dimethylethoxy)carbonyl]-, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester, [4aS-(4a.alpha.,6.alpha.,8aR*)]- (9CI)
MF C24 H32 N2 O6

Absolute stereochemistry.

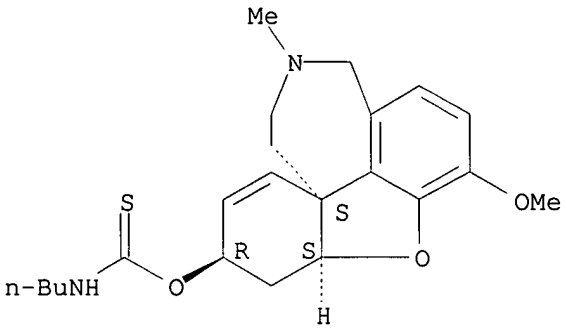


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamothioic acid, butyl-, O-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl) ester,

[4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C22 H30 N2 O3 S

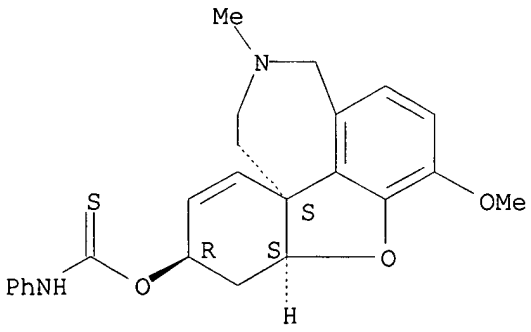
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamothioic acid, phenyl-, O-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl) ester, [4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C24 H26 N2 O3 S

Absolute stereochemistry.

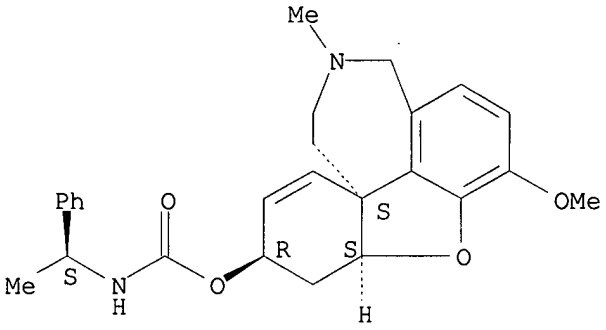


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN Carbamic acid, (1-phenylethyl)-, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester, [4aS-[4a.alpha.,6.beta.(R*),8aR*]]- (9CI)
MF C26 H30 N2 O4

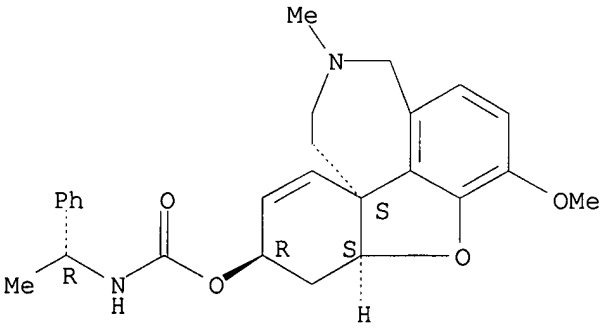
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, (1-phenylethyl)-, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester, [4aS-[4a.alpha.,6.beta.(S*),8aR*]]- (9CI)
MF C26 H30 N2 O4

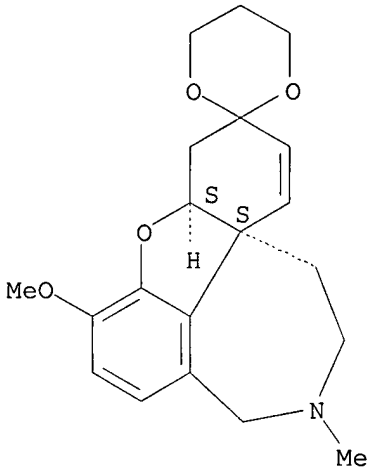
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxane], 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR*,8aR*)- (9CI)
MF C20 H25 N O4

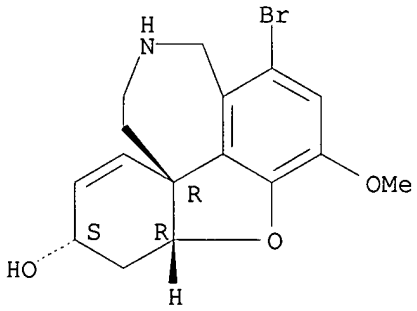
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-, [4aR-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C16 H18 Br N O3

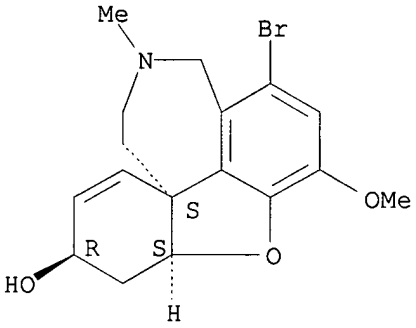
Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aS,6R,8aS)- (9CI)
MF C17 H20 Br N O3

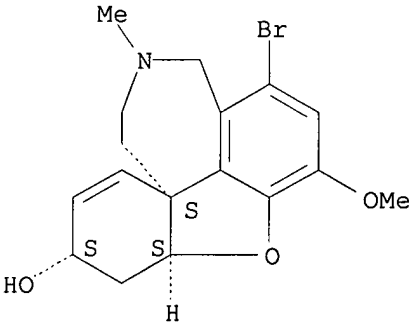
Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR,6R,8aR)-rel- (9CI)
MF C17 H20 Br N O3

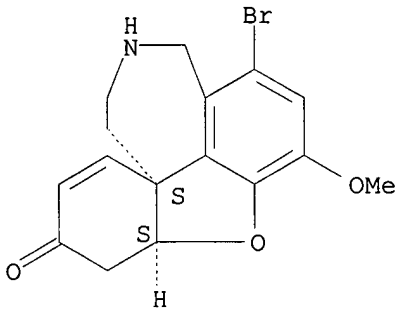
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4aR,8aR)-rel- (9CI)
MF C16 H16 Br N O3

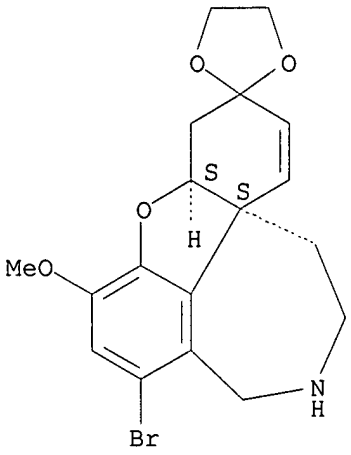
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxolane], 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4aR*,8aR*)- (9CI)
MF C18 H20 Br N O4

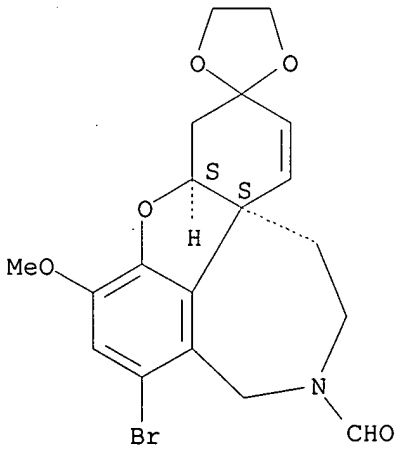
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxolane]-11(12H)-carboxaldehyde, 1-bromo-4a,5,9,10-tetrahydro-3-methoxy-, (4aR*,8aR*)-(9CI)
MF C19 H20 Br N O5

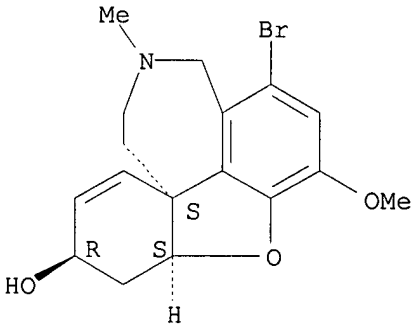
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR,6S,8aR)-rel- (9CI)
MF C17 H20 Br N O3

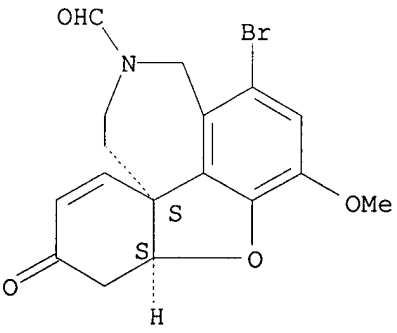
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

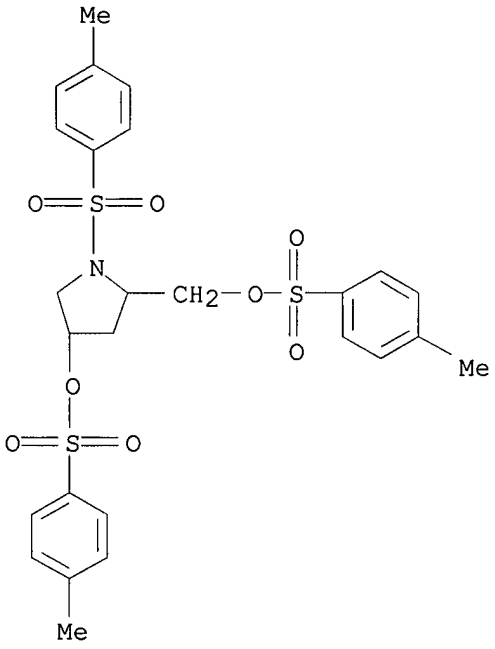
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-carboxaldehyde, 1-bromo-4a,5,9,10-tetrahydro-3-methoxy-6-oxo-, (4aR,8aR)-rel- (9CI)
MF C17 H16 Br N O4

Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

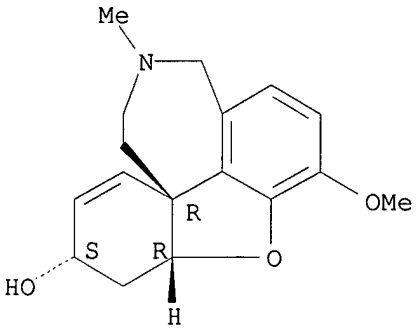
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2-Pyrrolidinemethanol, 1-[(4-methylphenyl)sulfonyl]-4-[[(4-methylphenyl)sulfonyl]oxy]-, 4-methylbenzenesulfonate (ester) (9CI)
MF C26 H29 N O8 S3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR,6S,8aR)- (9CI)
MF C17 H21 N O3
CI COM

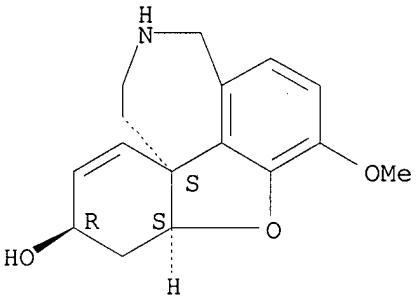
Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-, (4aR,6S,8aR)-rel- (9CI)
MF C16 H19 N O3

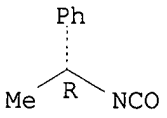
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzene, [(1R)-1-isocyanatoethyl]- (9CI)
MF C9 H9 N O
CI COM

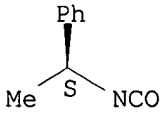
Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

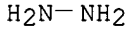
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzene, [(1S)-1-isocyanatoethyl]- (9CI)
MF C9 H9 N O

Absolute stereochemistry. Rotation (-).

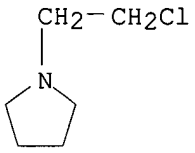


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazine, monohydrate (8CI, 9CI)
MF H4 N2 . H2 O
CI COM



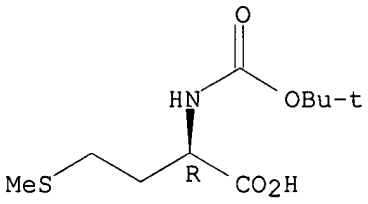
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Pyrrolidine, 1-(2-chloroethyl)-, hydrochloride (6CI, 7CI, 8CI, 9CI)
MF C6 H12 Cl N . Cl H



L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS

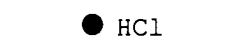
IN D-Methionine, N-[(1,1-dimethylethoxy)carbonyl]- (9CI)
MF C10 H19 N O4 S
CI COM

Absolute stereochemistry.

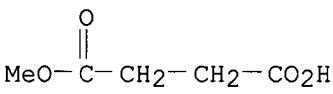


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Ethanamine, 2-chloro-N,N-dimethyl-, hydrochloride (9CI)
MF C4 H10 Cl N . Cl H



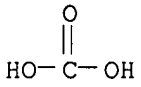
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Butanedioic acid, monomethyl ester (9CI)
MF C5 H8 O4
CI COM



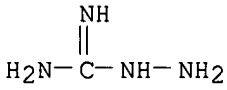
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbonic acid, compd. with hydrazinecarboximidamide (1:1) (9CI)
MF C H6 N4 . C H2 O3
CI COM

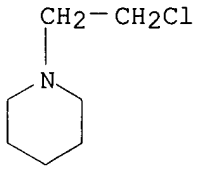
CM 1



CM 2

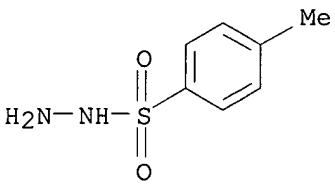


L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Piperidine, 1-(2-chloroethyl)-, hydrochloride (6CI, 7CI, 8CI, 9CI)
MF C7 H14 Cl N . Cl H



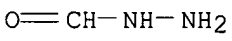
● HCl

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenesulfonic acid, 4-methyl-, hydrazide (9CI)
MF C7 H10 N2 O2 S
CI COM



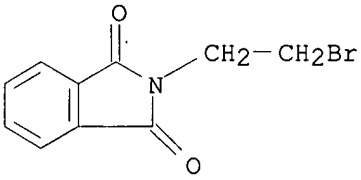
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazinecarboxaldehyde (9CI)
MF C H4 N2 O
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

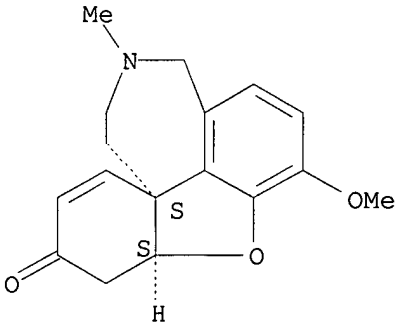
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 1H-Isoindole-1,3(2H)-dione, 2-(2-bromoethyl)- (9CI)
MF C10 H8 Br N O2

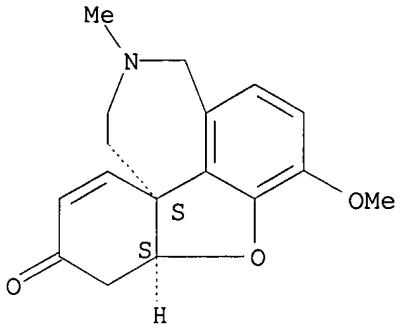


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aS,8aS)- (9CI)
MF C17 H19 N O3
CI COM

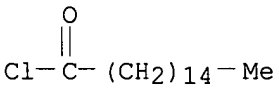
Absolute stereochemistry. Rotation (-).





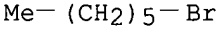
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hexadecanoyl chloride (9CI)
MF C16 H31 Cl O



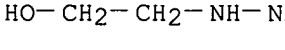
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hexane, 1-bromo- (6CI, 8CI, 9CI)
MF C6 H13 Br
CI COM



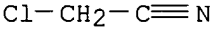
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Ethanol, 2-hydrazino- (6CI, 7CI, 8CI, 9CI)
MF C2 H8 N2 O
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Acetonitrile, chloro- (6CI, 8CI, 9CI)
MF C2 H2 Cl N
CI COM

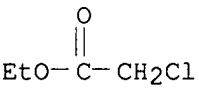


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 1-Propyne, 3-bromo- (9CI)
MF C3 H3 Br
CI COM

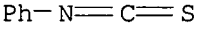
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

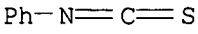
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Acetic acid, chloro-, ethyl ester (6CI, 8CI, 9CI)
MF C4 H7 Cl O2
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

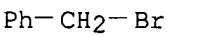
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzene, isothiocyanato- (9CI)
MF C7 H5 N S
CI COM





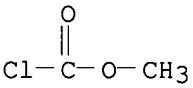
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzene, (bromomethyl)- (9CI)
MF C7 H7 Br
CI COM



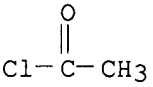
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbonochloridic acid, methyl ester (9CI)
MF C2 H3 Cl O2
CI COM



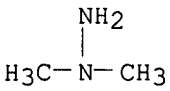
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Acetyl chloride (8CI, 9CI)
MF C2 H3 Cl O
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazine, 1,1-dimethyl- (8CI, 9CI)
MF C2 H8 N2
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> logoff
ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF
LOGOFF? (Y)/N/HOLD:y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.76	4.22

STN INTERNATIONAL LOGOFF AT 18:01:46 ON 04 OCT 2002